



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 22, 2021 – 10:07 AM EDT

PDB ID : 6WAI  
Title : Crystal Structure of SmcR N142D from *Vibrio vulnificus*  
Authors : Newman, J.D.; Russell, M.M.; Gonzalez-Gutierrez, G.; van Kessel, J.C.  
Deposited on : 2020-03-25  
Resolution : 2.58 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](mailto: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.

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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	:	2.20
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)	:	2.20

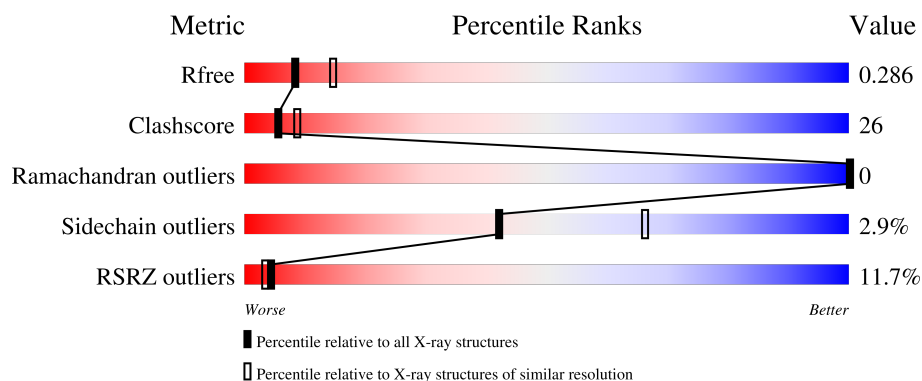
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.58 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	3676 (2.60-2.56)
Clashscore	141614	4049 (2.60-2.56)
Ramachandran outliers	138981	3979 (2.60-2.56)
Sidechain outliers	138945	3979 (2.60-2.56)
RSRZ outliers	127900	3614 (2.60-2.56)

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  $\geq 3$ , 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	225	<div> <div>13%</div> <div>51%</div> <div>36%</div> <div>•</div> <div>11%</div> </div>
1	B	225	<div> <div>10%</div> <div>48%</div> <div>38%</div> <div>•</div> <div>10%</div> </div>
1	C	225	<div> <div>9%</div> <div>58%</div> <div>32%</div> <div>•</div> <div>10%</div> </div>
1	D	225	<div> <div>10%</div> <div>48%</div> <div>38%</div> <div>•</div> <div>12%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6661 atoms, of which 12 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 LuxR family transcriptional regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	200	Total	C	N	O	S	0	0	0
			1617	1024	283	303	7			
1	B	202	Total	C	N	O	S	0	0	0
			1636	1035	289	305	7			
1	C	203	Total	C	N	O	S	0	0	0
			1648	1042	292	307	7			
1	D	199	Total	C	N	O	S	0	0	0
			1620	1025	289	299	7			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q9L8G8
A	-18	GLY	-	expression tag	UNP Q9L8G8
A	-17	SER	-	expression tag	UNP Q9L8G8
A	-16	SER	-	expression tag	UNP Q9L8G8
A	-15	HIS	-	expression tag	UNP Q9L8G8
A	-14	HIS	-	expression tag	UNP Q9L8G8
A	-13	HIS	-	expression tag	UNP Q9L8G8
A	-12	HIS	-	expression tag	UNP Q9L8G8
A	-11	HIS	-	expression tag	UNP Q9L8G8
A	-10	HIS	-	expression tag	UNP Q9L8G8
A	-9	SER	-	expression tag	UNP Q9L8G8
A	-8	SER	-	expression tag	UNP Q9L8G8
A	-7	GLY	-	expression tag	UNP Q9L8G8
A	-6	LEU	-	expression tag	UNP Q9L8G8
A	-5	VAL	-	expression tag	UNP Q9L8G8
A	-4	PRO	-	expression tag	UNP Q9L8G8
A	-3	ARG	-	expression tag	UNP Q9L8G8
A	-2	GLY	-	expression tag	UNP Q9L8G8
A	-1	SER	-	expression tag	UNP Q9L8G8
A	0	HIS	-	expression tag	UNP Q9L8G8
A	142	ASP	ASN	engineered mutation	UNP Q9L8G8

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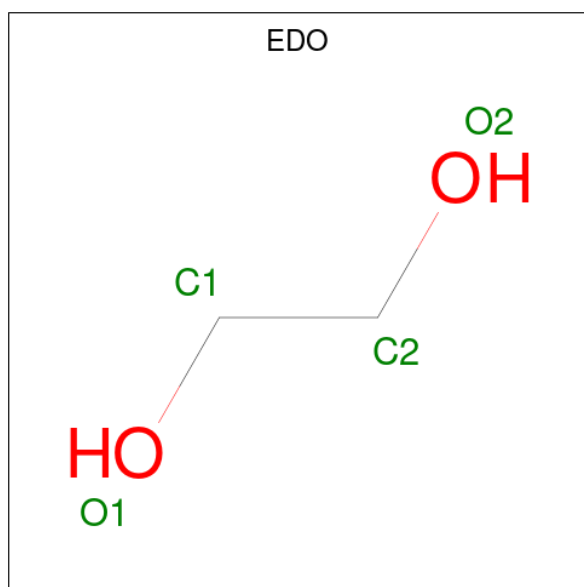
Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	MET	-	expression tag	UNP Q9L8G8
B	-18	GLY	-	expression tag	UNP Q9L8G8
B	-17	SER	-	expression tag	UNP Q9L8G8
B	-16	SER	-	expression tag	UNP Q9L8G8
B	-15	HIS	-	expression tag	UNP Q9L8G8
B	-14	HIS	-	expression tag	UNP Q9L8G8
B	-13	HIS	-	expression tag	UNP Q9L8G8
B	-12	HIS	-	expression tag	UNP Q9L8G8
B	-11	HIS	-	expression tag	UNP Q9L8G8
B	-10	HIS	-	expression tag	UNP Q9L8G8
B	-9	SER	-	expression tag	UNP Q9L8G8
B	-8	SER	-	expression tag	UNP Q9L8G8
B	-7	GLY	-	expression tag	UNP Q9L8G8
B	-6	LEU	-	expression tag	UNP Q9L8G8
B	-5	VAL	-	expression tag	UNP Q9L8G8
B	-4	PRO	-	expression tag	UNP Q9L8G8
B	-3	ARG	-	expression tag	UNP Q9L8G8
B	-2	GLY	-	expression tag	UNP Q9L8G8
B	-1	SER	-	expression tag	UNP Q9L8G8
B	0	HIS	-	expression tag	UNP Q9L8G8
B	142	ASP	ASN	engineered mutation	UNP Q9L8G8
C	-19	MET	-	expression tag	UNP Q9L8G8
C	-18	GLY	-	expression tag	UNP Q9L8G8
C	-17	SER	-	expression tag	UNP Q9L8G8
C	-16	SER	-	expression tag	UNP Q9L8G8
C	-15	HIS	-	expression tag	UNP Q9L8G8
C	-14	HIS	-	expression tag	UNP Q9L8G8
C	-13	HIS	-	expression tag	UNP Q9L8G8
C	-12	HIS	-	expression tag	UNP Q9L8G8
C	-11	HIS	-	expression tag	UNP Q9L8G8
C	-10	HIS	-	expression tag	UNP Q9L8G8
C	-9	SER	-	expression tag	UNP Q9L8G8
C	-8	SER	-	expression tag	UNP Q9L8G8
C	-7	GLY	-	expression tag	UNP Q9L8G8
C	-6	LEU	-	expression tag	UNP Q9L8G8
C	-5	VAL	-	expression tag	UNP Q9L8G8
C	-4	PRO	-	expression tag	UNP Q9L8G8
C	-3	ARG	-	expression tag	UNP Q9L8G8
C	-2	GLY	-	expression tag	UNP Q9L8G8
C	-1	SER	-	expression tag	UNP Q9L8G8
C	0	HIS	-	expression tag	UNP Q9L8G8
C	142	ASP	ASN	engineered mutation	UNP Q9L8G8

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-19	MET	-	expression tag	UNP Q9L8G8
D	-18	GLY	-	expression tag	UNP Q9L8G8
D	-17	SER	-	expression tag	UNP Q9L8G8
D	-16	SER	-	expression tag	UNP Q9L8G8
D	-15	HIS	-	expression tag	UNP Q9L8G8
D	-14	HIS	-	expression tag	UNP Q9L8G8
D	-13	HIS	-	expression tag	UNP Q9L8G8
D	-12	HIS	-	expression tag	UNP Q9L8G8
D	-11	HIS	-	expression tag	UNP Q9L8G8
D	-10	HIS	-	expression tag	UNP Q9L8G8
D	-9	SER	-	expression tag	UNP Q9L8G8
D	-8	SER	-	expression tag	UNP Q9L8G8
D	-7	GLY	-	expression tag	UNP Q9L8G8
D	-6	LEU	-	expression tag	UNP Q9L8G8
D	-5	VAL	-	expression tag	UNP Q9L8G8
D	-4	PRO	-	expression tag	UNP Q9L8G8
D	-3	ARG	-	expression tag	UNP Q9L8G8
D	-2	GLY	-	expression tag	UNP Q9L8G8
D	-1	SER	-	expression tag	UNP Q9L8G8
D	0	HIS	-	expression tag	UNP Q9L8G8
D	142	ASP	ASN	engineered mutation	UNP Q9L8G8

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	H	O	0	0
			10	2	6	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

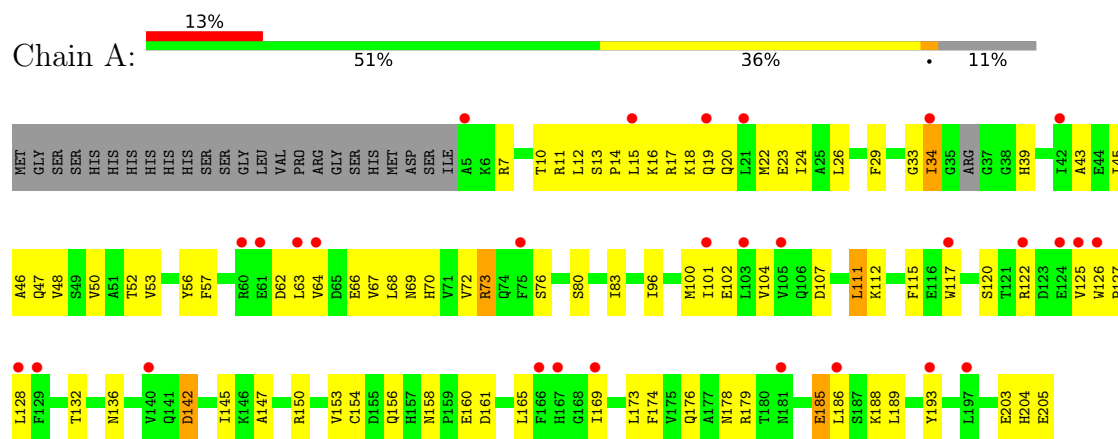
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	22	Total	O	0	0
			22	22		
4	B	29	Total	O	0	0
			29	29		
4	C	33	Total	O	0	0
			33	33		
4	D	31	Total	O	0	0
			31	31		

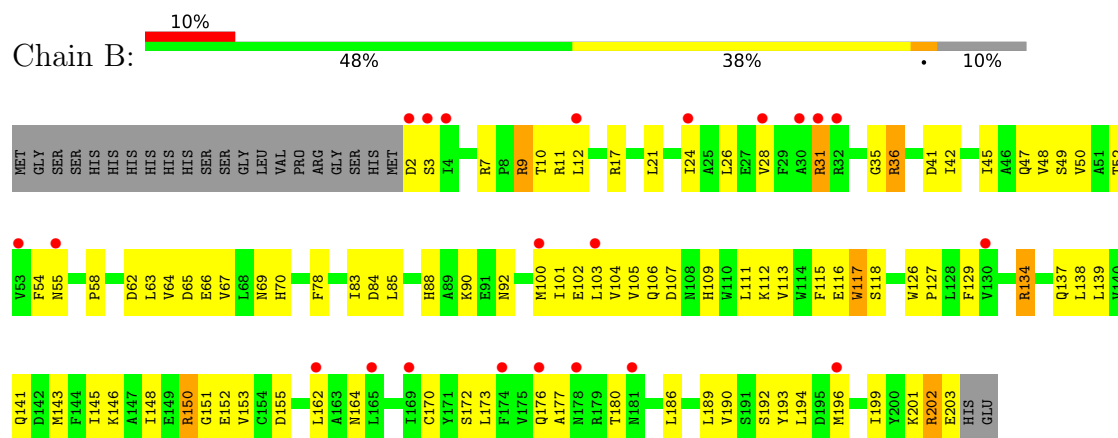
### 3 Residue-property plots [i](#)

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.

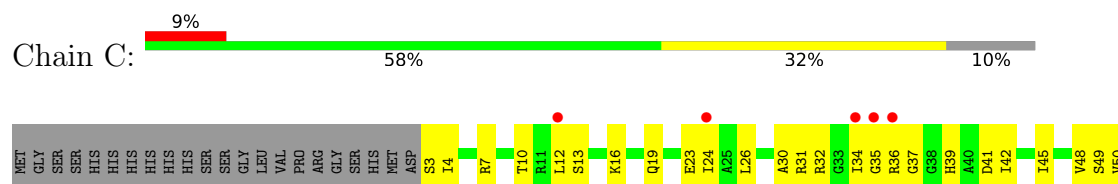
- Molecule 1: LuxR family transcriptional regulator

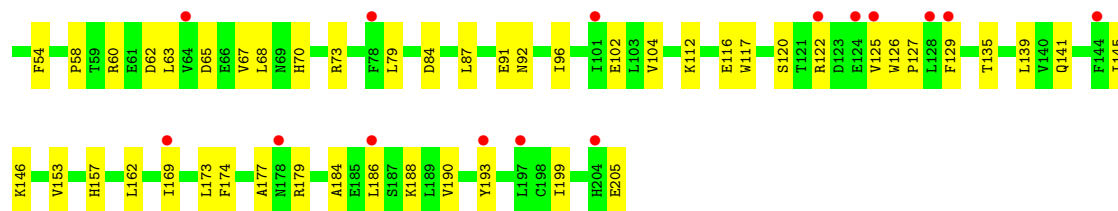


- Molecule 1: LuxR family transcriptional regulator

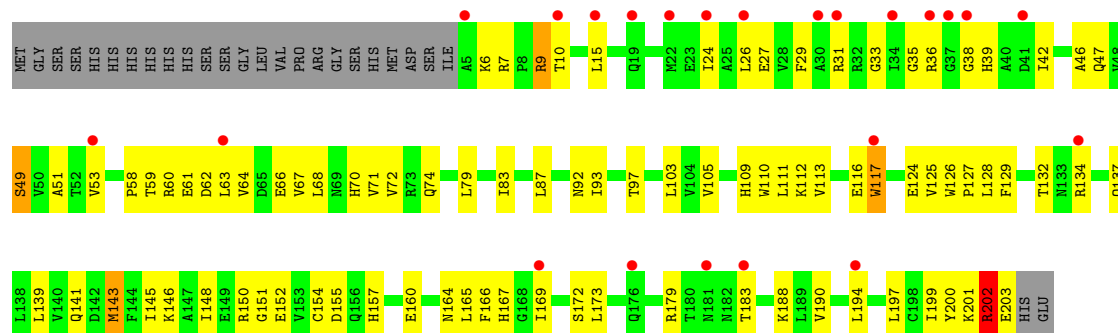


- Molecule 1: LuxR family transcriptional regulator





● Molecule 1: LuxR family transcriptional regulator





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	78.02Å 99.56Å 129.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	78.92 – 2.58 78.93 – 2.58	Depositor EDS
% Data completeness (in resolution range)	99.6 (78.92-2.58) 99.8 (78.93-2.58)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.17	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.90 (at 2.58Å)	Xtriage
Refinement program	PHENIX (1.15_3459: ???)	Depositor
R, $R_{free}$	0.231 , 0.286 0.231 , 0.286	Depositor DCC
$R_{free}$ test set	1670 reflections (5.18%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	42.4	Xtriage
Anisotropy	0.803	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 40.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	6661	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 36.73 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.7763e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, EDO

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	0.27	0/1649	0.44	0/2236
1	B	0.27	0/1668	0.49	1/2263 (0.0%)
1	C	0.26	0/1681	0.43	0/2279
1	D	0.28	0/1652	0.47	1/2240 (0.0%)
All	All	0.27	0/6650	0.46	2/9018 (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	A	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	202	ARG	NE-CZ-NH1	7.25	123.92	120.30
1	D	202	ARG	NE-CZ-NH2	-6.66	116.97	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	204	HIS	Peptide
1	A	34	ILE	Peptide

## 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	1617	0	1578	100	1
1	B	1636	0	1610	86	1
1	C	1648	0	1619	65	1
1	D	1620	0	1601	96	3
2	A	4	6	6	2	0
2	C	4	6	6	2	0
3	C	5	0	0	0	0
4	A	22	0	0	8	0
4	B	29	0	0	4	0
4	C	33	0	0	2	0
4	D	31	0	0	9	0
All	All	6649	12	6420	333	3

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 26.

All (333) 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:152:GLU:HG2	1:D:202:ARG:HG2	1.34	1.06
1:D:83:ILE:HD12	1:D:139:LEU:HD21	1.44	1.00
1:D:199:ILE:O	4:D:301:HOH:O	1.80	0.98
1:D:146:LYS:NZ	4:D:302:HOH:O	1.97	0.96
1:D:152:GLU:HG2	1:D:202:ARG:CG	1.93	0.96
1:D:83:ILE:HD12	1:D:139:LEU:CD2	1.97	0.93
1:D:167:HIS:ND1	4:D:303:HOH:O	2.00	0.92
1:B:137:GLN:NE2	4:B:301:HOH:O	2.05	0.88
1:D:9:ARG:NH1	1:D:10:THR:O	2.06	0.87
1:A:154:CYS:SG	4:A:406:HOH:O	2.37	0.83
1:A:73:ARG:NH2	4:A:402:HOH:O	2.06	0.82
1:B:69:ASN:ND2	4:B:302:HOH:O	2.12	0.82
1:C:39:HIS:HE1	1:C:60:ARG:HA	1.44	0.82
1:C:91:GLU:OE2	4:C:401:HOH:O	1.98	0.80
1:A:176:GLN:NE2	1:A:189:LEU:HD21	1.98	0.79
1:B:66:GLU:HG3	4:B:308:HOH:O	1.83	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:101:ILE:O	1:B:105:VAL:HG23	1.82	0.78
1:D:87:LEU:O	4:D:304:HOH:O	2.01	0.78
1:A:39:HIS:HB3	1:A:53:VAL:HG21	1.66	0.78
1:B:41:ASP:O	1:B:45:ILE:HD12	1.84	0.77
1:A:69:ASN:O	4:A:401:HOH:O	2.01	0.77
1:A:154:CYS:HB3	1:A:156:GLN:NE2	2.00	0.77
1:A:203:GLU:N	4:A:403:HOH:O	2.11	0.76
1:A:10:THR:O	1:A:17:ARG:NH2	2.17	0.76
1:B:151:GLY:C	1:B:202:ARG:HH21	1.89	0.75
1:D:202:ARG:HA	1:D:202:ARG:NE	2.01	0.74
1:B:164:ASN:ND2	4:B:303:HOH:O	2.15	0.74
1:C:13:SER:H	1:C:16:LYS:HE2	1.51	0.74
1:B:112:LYS:O	1:B:116:GLU:HG3	1.88	0.73
1:C:41:ASP:O	1:C:45:ILE:HD12	1.89	0.73
1:D:148:ILE:HG23	1:D:155:ASP:HA	1.71	0.73
1:A:24:ILE:HD13	1:A:45:ILE:HG22	1.71	0.72
2:A:301:EDO:H22	1:B:172:SER:HA	1.71	0.72
1:C:7:ARG:HG3	1:C:24:ILE:HD11	1.71	0.72
1:B:107:ASP:HA	1:B:112:LYS:HE3	1.73	0.71
1:D:83:ILE:CD1	1:D:139:LEU:CD2	2.68	0.71
1:B:90:LYS:HB2	1:B:199:ILE:HG21	1.72	0.71
1:A:169:ILE:HG23	1:A:193:TYR:HD1	1.56	0.71
1:D:39:HIS:CD2	1:D:53:VAL:HG12	2.25	0.70
1:B:151:GLY:O	1:B:202:ARG:NH2	2.25	0.69
1:B:24:ILE:O	1:B:28:VAL:HG23	1.93	0.69
1:D:68:LEU:HD23	1:D:128:LEU:HD23	1.74	0.68
1:A:22:MET:CE	1:A:66:GLU:HB3	2.23	0.68
2:C:301:EDO:H22	1:D:172:SER:HA	1.75	0.68
1:C:73:ARG:HH12	1:C:135:THR:HG21	1.57	0.68
1:D:141:GLN:OE1	1:D:160:GLU:HG3	1.93	0.68
1:B:146:LYS:O	1:B:150:ARG:HG2	1.93	0.68
1:A:57:PHE:CD1	1:A:63:LEU:HD13	2.28	0.67
1:A:73:ARG:HG2	4:A:401:HOH:O	1.95	0.67
1:C:122:ARG:HB2	1:C:125:VAL:CG2	2.24	0.67
1:B:100:MET:HE1	1:B:170:CYS:HB3	1.77	0.67
1:A:161:ASP:OD2	4:A:404:HOH:O	2.13	0.67
1:D:112:LYS:O	1:D:116:GLU:HG3	1.95	0.66
1:A:72:VAL:HG21	1:A:132:THR:HG22	1.75	0.66
1:D:39:HIS:HD2	1:D:53:VAL:HG12	1.61	0.66
1:D:92:ASN:ND2	1:D:143:MET:SD	2.69	0.66
1:A:205:GLU:O	4:A:405:HOH:O	2.14	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:70:HIS:O	1:D:74:GLN:HG2	1.96	0.65
1:C:186:LEU:O	1:C:190:VAL:HG23	1.97	0.65
1:D:200:TYR:HB3	1:D:202:ARG:NH2	2.12	0.64
1:D:9:ARG:HD3	1:D:10:THR:H	1.62	0.64
1:A:165:LEU:HD22	1:B:196:MET:HE2	1.80	0.64
1:A:19:GLN:HA	1:A:22:MET:HB3	1.79	0.64
1:C:126:TRP:CG	1:C:127:PRO:HD3	2.33	0.64
1:C:126:TRP:CE2	1:D:179:ARG:HG2	2.32	0.64
1:C:122:ARG:HB2	1:C:125:VAL:HG22	1.80	0.64
1:D:26:LEU:HD11	1:D:70:HIS:CE1	2.32	0.63
1:A:176:GLN:OE1	1:B:164:ASN:HB3	1.99	0.63
1:A:15:LEU:HA	1:A:18:LYS:CB	2.28	0.63
1:B:58:PRO:HD2	1:B:62:ASP:OD2	1.98	0.63
1:A:12:LEU:HD12	1:A:17:ARG:HA	1.80	0.63
1:A:20:GLN:HA	1:A:23:GLU:OE1	1.99	0.63
1:A:39:HIS:HB3	1:A:53:VAL:CG2	2.29	0.63
1:B:17:ARG:O	1:B:21:LEU:HG	1.97	0.63
1:D:190:VAL:O	1:D:194:LEU:HG	1.99	0.62
1:D:42:ILE:CD1	1:D:63:LEU:HD21	2.30	0.62
1:A:15:LEU:HA	1:A:18:LYS:HB3	1.81	0.61
1:C:188:LYS:HD3	1:C:188:LYS:N	2.15	0.61
1:D:68:LEU:CD2	1:D:128:LEU:HD23	2.30	0.61
1:C:30:ALA:HB1	1:C:112:LYS:HD2	1.83	0.61
1:D:93:ILE:HG23	1:D:166:PHE:HE1	1.66	0.61
1:A:154:CYS:HB3	1:A:156:GLN:CD	2.21	0.60
1:B:78:PHE:HZ	1:B:100:MET:HG3	1.66	0.60
1:A:68:LEU:HD13	1:A:117:TRP:HE3	1.66	0.60
1:A:165:LEU:HD22	1:B:196:MET:CE	2.31	0.60
1:D:152:GLU:CG	1:D:202:ARG:HG2	2.23	0.60
1:A:15:LEU:O	1:A:18:LYS:HB3	2.01	0.59
1:B:31:ARG:O	1:B:31:ARG:HD3	2.02	0.59
1:B:88:HIS:HB3	1:B:152:GLU:OE1	2.03	0.58
1:D:59:THR:HG22	1:D:61:GLU:OE2	2.03	0.58
1:A:104:VAL:HG23	1:A:111:LEU:CD2	2.34	0.58
1:B:118:SER:HA	1:B:129:PHE:CE2	2.38	0.58
1:D:9:ARG:HD3	1:D:10:THR:N	2.19	0.58
1:D:60:ARG:O	1:D:64:VAL:HG23	2.03	0.58
1:A:142:ASP:HA	1:A:145:ILE:HD12	1.85	0.58
1:A:22:MET:HE3	1:A:66:GLU:HB3	1.85	0.57
1:B:107:ASP:OD2	1:B:112:LYS:HE3	2.04	0.57
1:D:63:LEU:O	1:D:67:VAL:HG23	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:GLU:O	1:A:189:LEU:N	2.31	0.57
1:A:11:ARG:O	1:A:12:LEU:HD23	2.04	0.57
1:A:205:GLU:OE2	1:A:205:GLU:N	2.38	0.57
1:B:101:ILE:HD11	1:B:173:LEU:HD12	1.85	0.57
1:C:13:SER:H	1:C:16:LYS:CE	2.16	0.57
1:D:83:ILE:CD1	1:D:139:LEU:HD23	2.32	0.57
1:A:10:THR:HG22	1:A:11:ARG:O	2.05	0.57
1:C:112:LYS:HG3	1:C:174:PHE:HZ	1.70	0.56
1:D:164:ASN:OD1	4:D:305:HOH:O	2.18	0.56
1:A:111:LEU:HD21	1:A:115:PHE:CE2	2.40	0.56
1:C:12:LEU:O	4:C:402:HOH:O	2.18	0.56
1:A:15:LEU:CA	1:A:18:LYS:HB3	2.36	0.56
1:A:7:ARG:CZ	1:A:24:ILE:HG12	2.34	0.56
2:A:301:EDO:C2	1:B:172:SER:HA	2.35	0.56
1:B:35:GLY:C	1:B:36:ARG:HD2	2.26	0.56
1:C:42:ILE:HD11	1:C:63:LEU:HD21	1.87	0.56
1:D:167:HIS:CE1	4:D:303:HOH:O	2.52	0.56
1:D:169:ILE:HD13	1:D:197:LEU:HD11	1.88	0.56
1:A:24:ILE:HD12	1:A:46:ALA:HB2	1.88	0.55
1:C:79:LEU:HD13	1:C:139:LEU:HD12	1.86	0.55
1:A:76:SER:HB2	1:A:136:ASN:OD1	2.06	0.55
1:C:48:VAL:HG23	1:C:49:SER:O	2.07	0.55
1:D:62:ASP:O	1:D:66:GLU:HB2	2.06	0.55
1:D:105:VAL:HG21	1:D:183:THR:CG2	2.37	0.55
1:A:158:ASN:OD1	1:A:160:GLU:N	2.37	0.55
1:B:153:VAL:HG21	1:B:162:LEU:HD11	1.87	0.55
1:A:52:THR:O	1:A:56:TYR:HD2	1.90	0.55
1:A:68:LEU:HD13	1:A:117:TRP:CE3	2.42	0.55
1:A:107:ASP:OD1	1:A:112:LYS:HE3	2.07	0.55
1:B:78:PHE:CZ	1:B:100:MET:HG3	2.42	0.55
1:B:150:ARG:O	1:B:202:ARG:NE	2.36	0.55
1:C:157:HIS:HB2	1:C:162:LEU:HD11	1.89	0.55
1:A:34:ILE:HD11	1:A:64:VAL:HG22	1.89	0.54
1:C:19:GLN:O	1:C:23:GLU:HG3	2.06	0.54
1:C:31:ARG:NH1	1:C:32:ARG:HE	2.05	0.54
1:B:100:MET:HE1	1:B:115:PHE:HZ	1.73	0.54
1:D:83:ILE:HD12	1:D:139:LEU:HD23	1.84	0.54
1:B:64:VAL:HG23	1:B:65:ASP:N	2.22	0.54
1:B:83:ILE:CD1	1:B:139:LEU:HG	2.38	0.54
1:B:88:HIS:HB3	1:B:152:GLU:CD	2.27	0.54
1:D:27:GLU:OE2	1:D:31:ARG:NH1	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:39:HIS:CE1	1:C:60:ARG:HA	2.35	0.54
1:A:18:LYS:NZ	1:A:66:GLU:OE1	2.31	0.53
1:B:9:ARG:HB2	1:B:47:GLN:OE1	2.08	0.53
1:B:141:GLN:O	1:B:145:ILE:HG13	2.07	0.53
2:C:301:EDO:C2	1:D:172:SER:HA	2.36	0.53
1:B:31:ARG:HD3	1:B:31:ARG:C	2.29	0.53
1:B:52:THR:O	1:B:55:ASN:HB3	2.09	0.53
1:C:92:ASN:O	1:C:96:ILE:HG13	2.07	0.53
1:D:93:ILE:HG23	1:D:166:PHE:CE1	2.43	0.53
1:A:13:SER:HB2	1:A:16:LYS:H	1.74	0.53
1:A:126:TRP:CG	1:A:127:PRO:HD3	2.43	0.53
1:C:141:GLN:O	1:C:145:ILE:HG13	2.08	0.53
1:C:3:SER:O	1:C:3:SER:OG	2.27	0.53
1:D:72:VAL:HG11	1:D:132:THR:HG22	1.90	0.53
1:D:200:TYR:CB	1:D:202:ARG:NH2	2.72	0.53
1:D:58:PRO:HG2	1:D:62:ASP:OD2	2.09	0.52
1:D:117:TRP:CD1	1:D:129:PHE:HB2	2.44	0.52
1:B:201:LYS:O	1:B:202:ARG:NH1	2.43	0.52
1:D:38:GLY:O	1:D:42:ILE:HG13	2.10	0.52
1:A:26:LEU:CD2	1:A:67:VAL:HG13	2.39	0.52
1:A:33:GLY:HA3	1:B:36:ARG:HH12	1.74	0.52
1:B:117:TRP:CD1	1:B:129:PHE:HB2	2.44	0.52
1:B:186:LEU:O	1:B:190:VAL:HG23	2.09	0.52
1:D:71:VAL:HG12	1:D:110:TRP:HB2	1.92	0.52
1:A:154:CYS:HB3	1:A:156:GLN:HE22	1.74	0.52
1:A:111:LEU:HD21	1:A:115:PHE:HE2	1.74	0.52
1:D:103:LEU:HB2	1:D:111:LEU:HD23	1.92	0.52
1:D:202:ARG:NE	1:D:202:ARG:CA	2.72	0.52
1:A:104:VAL:HG23	1:A:111:LEU:HD22	1.92	0.52
1:B:192:SER:O	1:B:196:MET:HG3	2.10	0.52
1:A:22:MET:HE1	1:A:66:GLU:HB3	1.92	0.51
1:D:152:GLU:CG	1:D:202:ARG:CG	2.79	0.51
1:D:169:ILE:CD1	1:D:197:LEU:HD11	2.40	0.51
1:D:61:GLU:OE2	1:D:61:GLU:N	2.44	0.51
1:C:179:ARG:HG2	1:D:126:TRP:CE2	2.45	0.51
1:C:179:ARG:HG2	1:D:126:TRP:CD2	2.44	0.51
1:A:14:PRO:O	1:A:56:TYR:CE1	2.64	0.51
1:B:134:ARG:O	1:B:138:LEU:HG	2.10	0.51
1:B:172:SER:O	1:B:176:GLN:HG3	2.10	0.51
1:A:14:PRO:O	1:A:15:LEU:HD12	2.10	0.51
1:A:120:SER:OG	1:A:125:VAL:HG11	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:126:TRP:CG	1:B:127:PRO:HD3	2.46	0.51
1:D:7:ARG:HG3	1:D:24:ILE:HD11	1.93	0.51
1:B:50:VAL:HG12	1:B:54:PHE:HD2	1.75	0.51
1:C:58:PRO:HD2	1:C:62:ASP:OD1	2.11	0.50
1:D:151:GLY:O	1:D:202:ARG:NH1	2.44	0.50
1:A:205:GLU:O	1:A:205:GLU:HG2	2.11	0.50
1:A:101:ILE:HD11	1:A:173:LEU:HB3	1.94	0.50
1:C:26:LEU:CD2	1:C:67:VAL:HG13	2.42	0.50
1:A:13:SER:HB2	1:A:16:LYS:HB3	1.92	0.50
1:C:120:SER:OG	1:C:125:VAL:CG2	2.60	0.50
1:B:104:VAL:HG11	1:B:177:ALA:HB3	1.93	0.49
1:C:4:ILE:HD11	1:C:31:ARG:HD3	1.93	0.49
1:A:57:PHE:CG	1:A:63:LEU:HD13	2.46	0.49
1:B:35:GLY:O	1:B:36:ARG:HD2	2.13	0.49
1:D:141:GLN:O	1:D:145:ILE:HG13	2.12	0.49
1:C:4:ILE:CD1	1:C:31:ARG:HD3	2.42	0.49
1:C:126:TRP:N	1:C:127:PRO:CD	2.76	0.49
1:D:15:LEU:O	1:D:15:LEU:HD12	2.13	0.49
1:D:109:HIS:O	1:D:113:VAL:HG23	2.13	0.49
1:C:31:ARG:HH11	1:C:32:ARG:HE	1.59	0.48
1:B:12:LEU:HD12	1:B:17:ARG:NH1	2.29	0.48
1:A:122:ARG:HB2	1:A:125:VAL:HG12	1.95	0.48
1:D:6:LYS:HD3	1:D:47:GLN:OE1	2.14	0.48
1:A:50:VAL:O	1:A:53:VAL:HG22	2.13	0.48
1:C:36:ARG:HH21	1:D:36:ARG:NH1	2.12	0.48
1:D:35:GLY:O	1:D:36:ARG:HG3	2.14	0.48
1:A:11:ARG:C	1:A:12:LEU:HD23	2.34	0.48
1:A:34:ILE:HG23	1:A:34:ILE:O	2.14	0.48
1:B:26:LEU:HD11	1:B:70:HIS:CE1	2.49	0.48
1:A:73:ARG:HD3	1:A:73:ARG:HA	1.65	0.47
1:D:24:ILE:HG22	1:D:42:ILE:HG23	1.96	0.47
1:C:42:ILE:CD1	1:C:63:LEU:HD21	2.43	0.47
1:A:165:LEU:HB2	1:B:196:MET:HE3	1.97	0.47
1:C:4:ILE:HD12	1:C:32:ARG:NH2	2.30	0.47
1:D:126:TRP:CG	1:D:127:PRO:HD3	2.50	0.47
1:B:193:TYR:HA	1:B:196:MET:CE	2.45	0.47
1:A:13:SER:CB	1:A:16:LYS:HB3	2.44	0.47
1:A:52:THR:O	1:A:56:TYR:CD2	2.67	0.47
1:B:126:TRP:N	1:B:127:PRO:CD	2.78	0.47
1:C:35:GLY:O	1:C:37:GLY:N	2.41	0.47
1:C:193:TYR:CE2	1:D:165:LEU:HB2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:49:SER:OG	1:D:51:ALA:N	2.48	0.47
1:A:43:ALA:O	1:A:47:GLN:N	2.48	0.47
1:B:63:LEU:O	1:B:67:VAL:HG23	2.15	0.47
1:C:104:VAL:HG21	1:C:174:PHE:HD1	1.80	0.47
1:D:203:GLU:OE2	4:D:306:HOH:O	2.20	0.47
1:A:34:ILE:HD12	1:A:117:TRP:CZ2	2.50	0.47
1:C:26:LEU:HD23	1:C:67:VAL:HG13	1.97	0.47
1:A:26:LEU:HD11	1:A:70:HIS:ND1	2.30	0.46
1:A:186:LEU:HD23	1:A:186:LEU:HA	1.78	0.46
1:D:26:LEU:CD2	1:D:67:VAL:HG13	2.45	0.46
1:A:57:PHE:CE1	1:A:63:LEU:HD13	2.50	0.46
1:C:16:LYS:HE2	1:C:16:LYS:HB2	1.74	0.46
1:A:179:ARG:HG2	1:B:126:TRP:CE2	2.51	0.46
1:A:100:MET:SD	1:A:101:ILE:HG12	2.56	0.46
1:A:43:ALA:HB1	1:A:48:VAL:O	2.15	0.46
1:A:126:TRP:N	1:A:127:PRO:CD	2.79	0.46
1:B:148:ILE:HG23	1:B:155:ASP:HA	1.98	0.46
1:B:176:GLN:O	1:B:180:THR:HG22	2.15	0.46
1:A:185:GLU:HA	1:A:188:LYS:HB3	1.98	0.45
1:A:64:VAL:O	1:A:68:LEU:HB2	2.16	0.45
1:D:68:LEU:HB2	1:D:117:TRP:CH2	2.51	0.45
1:A:19:GLN:O	1:A:23:GLU:N	2.44	0.45
1:A:72:VAL:HG21	1:A:132:THR:CG2	2.46	0.45
1:D:154:CYS:HB2	4:D:313:HOH:O	2.15	0.45
1:A:122:ARG:O	1:A:125:VAL:HG12	2.17	0.45
1:C:104:VAL:CG2	1:C:174:PHE:HD1	2.30	0.45
1:D:134:ARG:O	1:D:137:GLN:HB3	2.16	0.45
1:B:151:GLY:C	1:B:202:ARG:NH2	2.62	0.45
1:A:10:THR:HG22	1:A:12:LEU:HD23	1.97	0.45
1:C:10:THR:HG23	1:C:12:LEU:HG	1.99	0.45
1:D:103:LEU:CB	1:D:111:LEU:HD23	2.47	0.45
1:A:15:LEU:HA	1:A:18:LYS:HB2	1.99	0.45
1:B:92:ASN:HB3	1:B:143:MET:CE	2.46	0.45
1:A:147:ALA:HB1	1:A:153:VAL:HG23	1.98	0.44
1:D:124:GLU:HG2	1:D:125:VAL:HG23	2.00	0.44
1:D:126:TRP:N	1:D:127:PRO:CD	2.80	0.44
1:D:112:LYS:HE3	1:D:112:LYS:HB3	1.60	0.44
1:D:154:CYS:HB3	1:D:157:HIS:CD2	2.53	0.44
1:B:150:ARG:HB2	1:B:202:ARG:HG3	2.00	0.44
1:C:104:VAL:HG11	1:C:177:ALA:HB3	2.00	0.44
1:B:24:ILE:CG2	1:B:42:ILE:HG23	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:26:LEU:CD2	1:B:67:VAL:HG13	2.47	0.44
1:D:59:THR:HG22	1:D:60:ARG:N	2.33	0.44
1:A:29:PHE:CE2	1:A:34:ILE:HG13	2.52	0.44
1:C:184:ALA:O	1:C:188:LYS:HD3	2.18	0.44
1:B:10:THR:OG1	1:B:11:ARG:N	2.51	0.43
1:B:193:TYR:CD2	1:B:196:MET:HE1	2.53	0.43
1:B:102:GLU:O	1:B:106:GLN:HG3	2.17	0.43
1:B:134:ARG:NH1	1:B:138:LEU:HD21	2.33	0.43
1:C:36:ARG:HE	1:D:36:ARG:NH1	2.16	0.43
1:D:97:THR:CG2	1:D:173:LEU:HD13	2.49	0.43
1:B:84:ASP:O	1:B:92:ASN:ND2	2.42	0.43
1:D:79:LEU:HD13	1:D:139:LEU:HD22	1.99	0.43
1:D:202:ARG:HA	1:D:202:ARG:HE	1.81	0.43
1:A:15:LEU:HG	1:A:18:LYS:CB	2.48	0.43
1:A:15:LEU:C	1:A:18:LYS:HB3	2.38	0.43
1:A:128:LEU:N	1:A:128:LEU:HD22	2.34	0.43
1:B:137:GLN:O	1:B:141:GLN:HG3	2.19	0.43
1:B:152:GLU:HG3	1:B:202:ARG:HG2	2.01	0.43
1:B:193:TYR:HA	1:B:196:MET:HE2	2.01	0.43
1:C:122:ARG:CB	1:C:125:VAL:HG22	2.47	0.43
1:D:194:LEU:N	1:D:194:LEU:HD23	2.32	0.43
1:C:26:LEU:HD11	1:C:70:HIS:CE1	2.54	0.43
1:C:153:VAL:HG13	1:C:199:ILE:O	2.18	0.43
1:C:188:LYS:HD3	1:C:188:LYS:H	1.82	0.43
1:D:103:LEU:HB2	1:D:111:LEU:CD2	2.48	0.43
1:A:52:THR:HG23	1:A:56:TYR:HE2	1.83	0.43
1:B:83:ILE:HG12	1:B:143:MET:HE3	2.01	0.43
1:D:68:LEU:HD23	1:D:128:LEU:CD2	2.44	0.43
1:B:64:VAL:CG2	1:B:65:ASP:N	2.82	0.43
1:C:120:SER:OG	1:C:125:VAL:HG21	2.19	0.43
1:D:154:CYS:CA	4:D:313:HOH:O	2.67	0.42
1:A:128:LEU:HD13	1:A:128:LEU:HA	1.75	0.42
1:C:173:LEU:HG	1:C:193:TYR:CD1	2.54	0.42
1:D:201:LYS:C	1:D:202:ARG:HE	2.22	0.42
1:B:50:VAL:HG12	1:B:54:PHE:CD2	2.53	0.42
1:D:97:THR:HG21	1:D:173:LEU:HD13	2.00	0.42
1:C:102:GLU:OE1	1:C:102:GLU:HA	2.19	0.42
1:A:26:LEU:HD11	1:A:70:HIS:CE1	2.55	0.42
1:C:84:ASP:HB3	1:C:87:LEU:HG	2.02	0.42
1:C:126:TRP:CD2	1:C:127:PRO:HD3	2.55	0.42
1:D:26:LEU:HD11	1:D:70:HIS:ND1	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:GLN:O	1:A:22:MET:N	2.52	0.42
1:A:83:ILE:HG12	1:A:96:ILE:HD11	2.02	0.42
1:C:117:TRP:CD1	1:C:129:PHE:HB2	2.54	0.42
1:B:203:GLU:OE1	1:B:203:GLU:N	2.46	0.42
1:A:174:PHE:O	1:A:178:ASN:ND2	2.53	0.41
1:B:189:LEU:HD11	1:B:193:TYR:CE2	2.55	0.41
1:C:120:SER:OG	1:C:125:VAL:HG23	2.19	0.41
1:C:169:ILE:HG23	1:C:193:TYR:HD1	1.85	0.41
1:A:176:GLN:OE1	1:B:164:ASN:CB	2.66	0.41
1:C:68:LEU:HB2	1:C:117:TRP:CH2	2.55	0.41
1:D:148:ILE:HD13	1:D:157:HIS:O	2.20	0.41
1:A:185:GLU:CD	1:A:188:LYS:HD3	2.41	0.41
1:B:48:VAL:HG23	1:B:49:SER:O	2.20	0.41
1:C:146:LYS:HD2	1:C:146:LYS:HA	1.90	0.41
1:D:29:PHE:O	1:D:33:GLY:N	2.52	0.41
1:A:165:LEU:HD13	1:B:193:TYR:CD1	2.55	0.41
1:B:90:LYS:HE3	1:B:194:LEU:HB3	2.02	0.41
1:B:109:HIS:O	1:B:113:VAL:HG23	2.21	0.41
1:C:13:SER:N	1:C:16:LYS:HE2	2.26	0.41
1:D:173:LEU:HD13	1:D:190:VAL:HG22	2.02	0.41
1:D:202:ARG:CA	1:D:202:ARG:HE	2.32	0.41
1:A:150:ARG:NH1	4:A:408:HOH:O	2.47	0.41
1:C:34:ILE:HD11	1:C:116:GLU:HB2	2.03	0.41
1:D:150:ARG:O	1:D:202:ARG:HD2	2.21	0.41
1:B:100:MET:HE1	1:B:115:PHE:CZ	2.55	0.40
1:C:50:VAL:HG12	1:C:54:PHE:CD2	2.56	0.40
1:D:24:ILE:HD13	1:D:46:ALA:HB2	2.03	0.40
1:B:189:LEU:HA	1:B:189:LEU:HD12	1.74	0.40
1:B:85:LEU:HD22	1:B:146:LYS:HD2	2.03	0.40
1:B:103:LEU:HD12	1:B:111:LEU:HD22	2.03	0.40

All (3) 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:62:ASP:OD1	1:D:188:LYS:NZ[4_545]	2.13	0.07
1:B:55:ASN:O	1:D:134:ARG:NH1[3_654]	2.14	0.06
1:C:205:GLU:O	1:D:9:ARG:NH2[1_455]	2.16	0.04

## 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	196/225 (87%)	193 (98%)	3 (2%)	0	100	100
1	B	200/225 (89%)	198 (99%)	2 (1%)	0	100	100
1	C	201/225 (89%)	195 (97%)	6 (3%)	0	100	100
1	D	197/225 (88%)	193 (98%)	4 (2%)	0	100	100
All	All	794/900 (88%)	779 (98%)	15 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	177/201 (88%)	171 (97%)	6 (3%)	37	60
1	B	180/201 (90%)	171 (95%)	9 (5%)	24	45
1	C	181/201 (90%)	180 (99%)	1 (1%)	86	94
1	D	178/201 (89%)	173 (97%)	5 (3%)	43	67
All	All	716/804 (89%)	695 (97%)	21 (3%)	42	66

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	ARG
1	A	80	SER

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Mol	Chain	Res	Type
1	A	102	GLU
1	A	111	LEU
1	A	142	ASP
1	A	185	GLU
1	B	2	ASP
1	B	3	SER
1	B	7	ARG
1	B	9	ARG
1	B	31	ARG
1	B	36	ARG
1	B	117	TRP
1	B	134	ARG
1	B	150	ARG
1	C	65	ASP
1	D	9	ARG
1	D	49	SER
1	D	117	TRP
1	D	143	MET
1	D	202	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	164	ASN
1	C	156	GLN
1	D	164	ASN

### 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 ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry

3 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	EDO	A	301	-	3,3,3	0.46	0	2,2,2	0.34	0
3	SO4	C	302	-	4,4,4	0.15	0	6,6,6	0.06	0
2	EDO	C	301	-	3,3,3	0.47	0	2,2,2	0.31	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
2	EDO	A	301	-	-	1/1/1/1	-
2	EDO	C	301	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	301	EDO	O1-C1-C2-O2
2	A	301	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	EDO	2	0
2	C	301	EDO	2	0

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	200/225 (88%)	0.92	29 (14%) <b>2</b> <b>1</b>	35, 58, 93, 105	0
1	B	202/225 (89%)	0.78	22 (10%) <b>5</b> <b>4</b>	32, 51, 74, 92	0
1	C	203/225 (90%)	0.73	20 (9%) <b>7</b> <b>6</b>	28, 49, 75, 113	0
1	D	199/225 (88%)	0.84	23 (11%) <b>4</b> <b>3</b>	31, 55, 91, 117	0
All	All	804/900 (89%)	0.81	94 (11%) <b>4</b> <b>3</b>	28, 53, 88, 117	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	37	GLY	10.1
1	A	64	VAL	7.1
1	D	5	ALA	6.4
1	A	60	ARG	5.9
1	C	34	ILE	5.5
1	C	35	GLY	5.3
1	B	4	ILE	5.1
1	A	125	VAL	5.0
1	A	63	LEU	4.4
1	A	124	GLU	4.4
1	D	34	ILE	4.3
1	C	204	HIS	4.1
1	D	30	ALA	4.0
1	A	21	LEU	4.0
1	A	117	TRP	3.9
1	B	31	ARG	3.9
1	D	24	ILE	3.9
1	C	36	ARG	3.9
1	B	2	ASP	3.8
1	D	36	ARG	3.7
1	B	165	LEU	3.6

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Mol	Chain	Res	Type	RSRZ
1	D	53	VAL	3.6
1	D	15	LEU	3.3
1	D	38	GLY	3.2
1	A	34	ILE	3.2
1	B	32	ARG	3.2
1	C	128	LEU	3.2
1	B	169	ILE	3.2
1	C	124	GLU	3.2
1	C	144	PHE	3.1
1	D	169	ILE	3.1
1	D	181	ASN	3.1
1	A	122	ARG	3.0
1	A	169	ILE	2.9
1	A	5	ALA	2.9
1	A	129	PHE	2.9
1	C	197	LEU	2.9
1	B	3	SER	2.8
1	C	101	ILE	2.8
1	B	30	ALA	2.8
1	A	105	VAL	2.8
1	B	176	GLN	2.8
1	A	103	LEU	2.8
1	D	63	LEU	2.8
1	D	117	TRP	2.8
1	D	26	LEU	2.7
1	A	128	LEU	2.7
1	B	100	MET	2.7
1	C	129	PHE	2.7
1	A	15	LEU	2.7
1	D	134	ARG	2.7
1	A	140	VAL	2.6
1	D	22	MET	2.6
1	B	28	VAL	2.5
1	A	42	ILE	2.5
1	A	181	ASN	2.5
1	C	186	LEU	2.5
1	A	101	ILE	2.5
1	B	181	ASN	2.5
1	A	19	GLN	2.5
1	A	61	GLU	2.5
1	B	55	ASN	2.5
1	B	24	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	31	ARG	2.4
1	B	12	LEU	2.4
1	B	53	VAL	2.4
1	D	194	LEU	2.4
1	A	166	PHE	2.4
1	B	130	VAL	2.3
1	C	64	VAL	2.3
1	B	196	MET	2.3
1	C	78	PHE	2.3
1	C	178	ASN	2.3
1	C	12	LEU	2.3
1	B	174	PHE	2.3
1	B	103	LEU	2.3
1	C	169	ILE	2.2
1	A	167	HIS	2.2
1	C	193	TYR	2.2
1	D	10	THR	2.2
1	B	178	ASN	2.2
1	A	193	TYR	2.2
1	C	122	ARG	2.2
1	B	162	LEU	2.2
1	D	19	GLN	2.1
1	A	126	TRP	2.1
1	D	183	THR	2.1
1	D	176	GLN	2.1
1	C	125	VAL	2.1
1	A	75	PHE	2.1
1	A	186	LEU	2.0
1	C	24	ILE	2.0
1	A	197	LEU	2.0
1	D	41	ASP	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

## 6.3 Carbohydrates [i](#)

There are no 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, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	EDO	A	301	4/4	0.83	0.53	67,81,82,82	0
2	EDO	C	301	4/4	0.94	0.22	46,55,56,56	0
3	SO4	C	302	5/5	0.95	0.33	73,73,74,75	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.