



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 02:17 am GMT

PDB ID : 1O79  
Title : Structures of human oxidosqualene cyclase inhibitors bound to a homologous enzyme  
Authors : Lenhart, A.; Reinert, D.J.; Weihofen, W.A.; Aebi, J.D.; Dehmlow, H.; Morand, O.H.; Schulz, G.E.  
Deposited on : 2002-10-27  
Resolution : 2.80 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

<http://wwpdb.org/validation/2016/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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

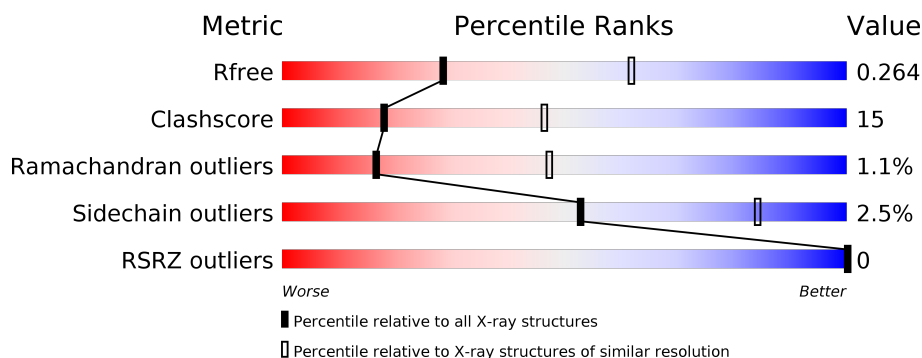
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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))
$R_{free}$	100719	2583 (2.80-2.80)
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)
RSRZ outliers	101464	2610 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	631	<div> <div>65%</div> <div>32%</div> <div>..</div> </div>
1	B	631	<div> <div>67%</div> <div>30%</div> <div>..</div> </div>
1	C	631	<div> <div>65%</div> <div>32%</div> <div>..</div> </div>

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	C8E	A	700	-	-	-	X
2	C8E	B	700	-	-	-	X
2	C8E	C	700	-	-	-	X
3	R23	A	800	-	-	-	X

## 2 Entry composition [i](#)

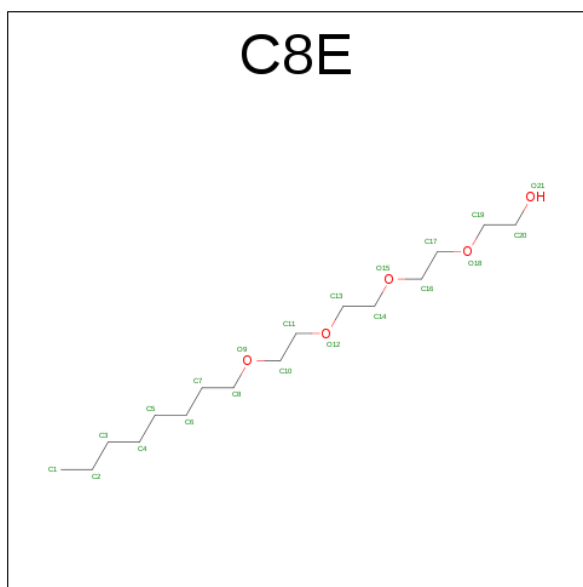
There are 4 unique types of molecules in this entry. The entry contains 15285 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SQUALENE-HOPENE CYCLASE.

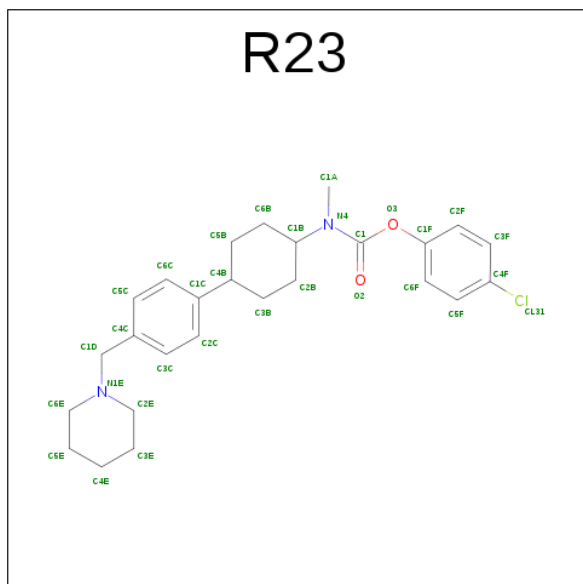
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	620	Total	C	N	O	S	0	0	1
			4964	3188	859	897	20			
1	B	620	Total	C	N	O	S	0	0	1
			4964	3188	859	897	20			
1	C	620	Total	C	N	O	S	0	0	1
			4964	3188	859	897	20			

- Molecule 2 is (HYDROXYETHYLOXY)TRI(ETHYLOXY)OCTANE (three-letter code: C8E) (formula:  $C_{16}H_{34}O_5$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			21	16	5		
2	B	1	Total	C	O	0	0
			21	16	5		
2	C	1	Total	C	O	0	0
			21	16	5		

- Molecule 3 is METHYL-[4-(4-PIPERIDINE-1-YLMETHYL-PHENYL)-CYCLOHEXYL]-CARBAMINIC ACID-(4-CHLOROPHENYL)-ESTER (three-letter code: R23) (formula:  $C_{26}H_{33}ClN_2O_2$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	0	0
			31	26	1	2	2		
3	B	1	Total	C	Cl	N	O	0	0
			31	26	1	2	2		
3	C	1	Total	C	Cl	N	O	0	0
			31	26	1	2	2		

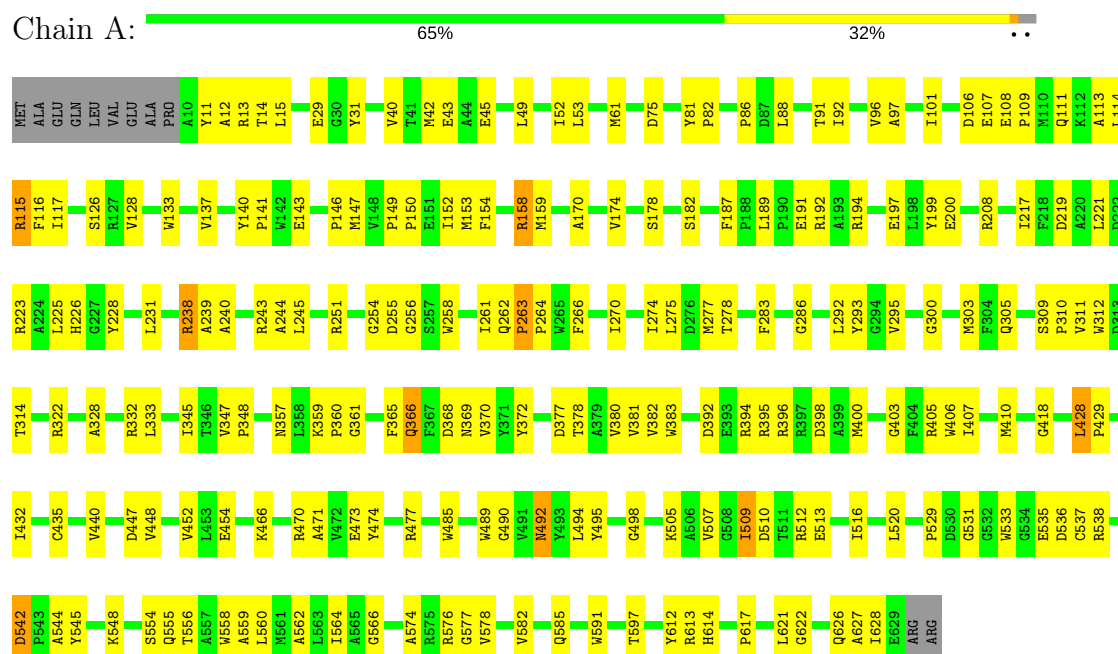
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	78	Total	O	0	0
			78	78		
4	B	84	Total	O	0	0
			84	84		
4	C	75	Total	O	0	0
			75	75		

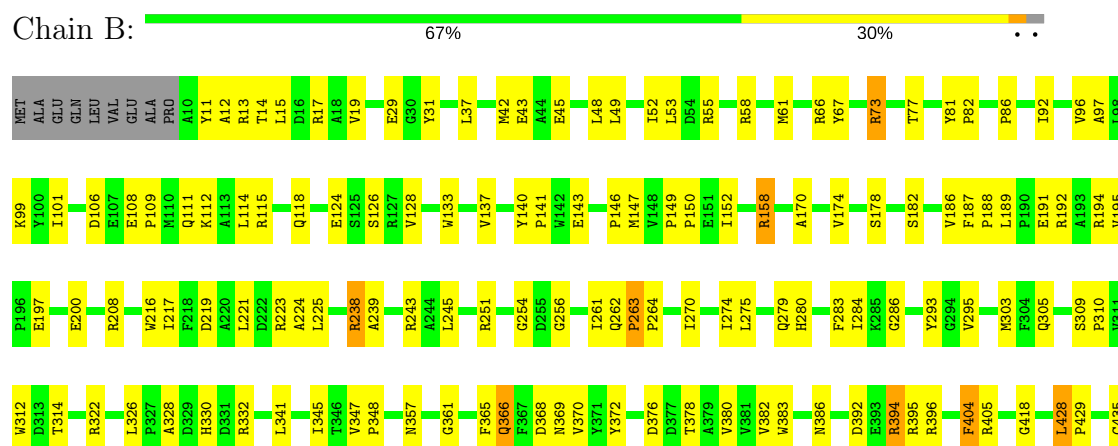
### 3 Residue-property plots

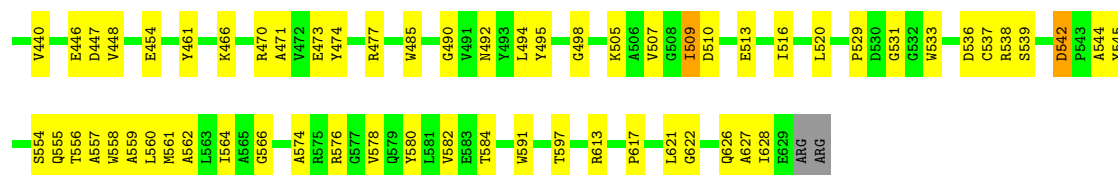
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: SQUALENE-HOPENE CYCLASE



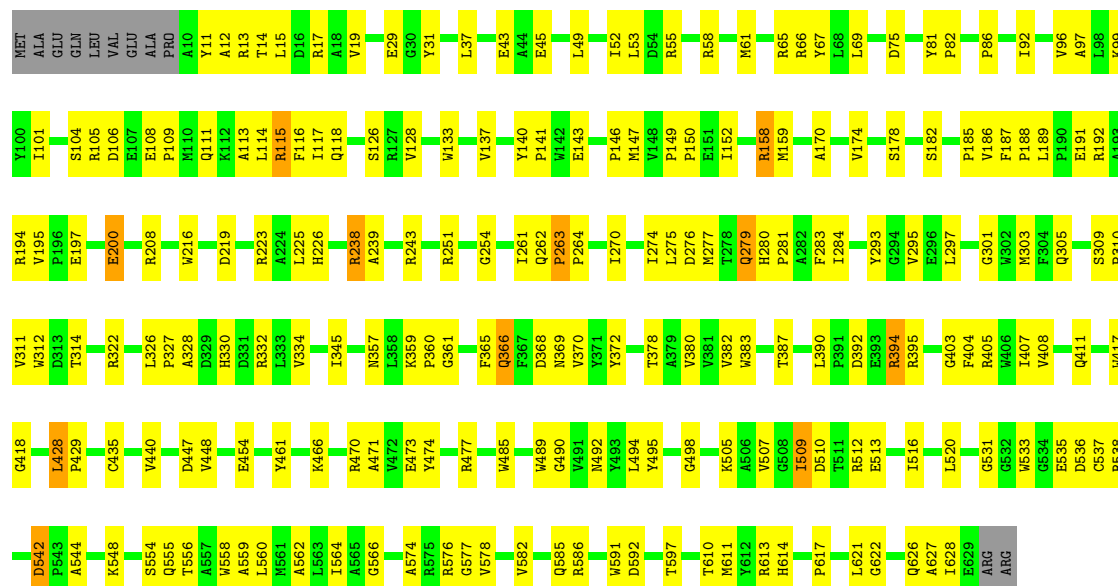
#### • Molecule 1: SQUALENE-HOPENE CYCLASE





• Molecule 1: SQUALENE-HOPENE CYCLASE

Chain C: 65% 32%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	141.29Å 141.29Å 245.40Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.00 – 2.80 32.43 – 2.80	Depositor EDS
% Data completeness (in resolution range)	74.2 (25.00-2.80) 74.2 (32.43-2.80)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.70 (at 2.81Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.239 , 0.273 0.235 , 0.264	Depositor DCC
$R_{free}$ test set	1979 reflections (3.94%)	DCC
Wilson B-factor (Å <sup>2</sup> )	25.5	Xtriage
Anisotropy	1.051	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.26 , 21.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.41$ , $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.008 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	15285	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	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 24.91 % 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 3.5044e-03. 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: R23, C8E

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	0.30	0/5117	0.52	0/6967
1	B	0.31	0/5117	0.53	0/6967
1	C	0.30	0/5117	0.52	0/6967
All	All	0.30	0/15351	0.52	0/20901

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	4964	0	4787	150	0
1	B	4964	0	4787	142	0
1	C	4964	0	4787	157	0
2	A	21	0	34	1	0
2	B	21	0	34	0	0
2	C	21	0	34	0	0
3	A	31	0	33	0	0
3	B	31	0	33	2	0
3	C	31	0	33	0	0
4	A	78	0	0	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	84	0	0	7	0
4	C	75	0	0	3	0
All	All	15285	0	14562	447	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 15.

All (447) 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:238:ARG:HB2	1:C:238:ARG:HH11	1.21	1.03
1:C:520:LEU:HD21	1:C:566:GLY:HA3	1.43	0.98
1:B:520:LEU:HD21	1:B:566:GLY:HA3	1.44	0.98
1:A:520:LEU:HD21	1:A:566:GLY:HA3	1.42	0.97
1:A:238:ARG:HB2	1:A:238:ARG:HH11	1.36	0.90
1:B:217:ILE:H	1:B:217:ILE:HD12	1.37	0.87
1:A:217:ILE:HD12	1:A:217:ILE:H	1.44	0.83
1:A:239:ALA:O	1:A:243:ARG:HG2	1.82	0.79
1:A:489:TRP:HB2	4:A:2069:HOH:O	1.88	0.73
1:C:97:ALA:O	1:C:101:ILE:HG13	1.89	0.73
1:B:97:ALA:O	1:B:101:ILE:HG13	1.88	0.72
1:A:277:MET:HG2	4:A:2041:HOH:O	1.88	0.72
1:C:627:ALA:C	1:C:628:ILE:HD12	2.11	0.71
1:A:544:ALA:O	1:A:548:LYS:HD3	1.90	0.70
1:C:309:SER:HB3	1:C:365:PHE:CZ	2.26	0.70
1:A:159:MET:HG3	4:A:2027:HOH:O	1.91	0.70
1:C:159:MET:HG3	4:C:2025:HOH:O	1.91	0.70
1:C:622:GLY:O	1:C:626:GLN:HG2	1.93	0.69
1:B:170:ALA:O	1:B:174:VAL:HG23	1.91	0.69
1:A:309:SER:HB3	1:A:365:PHE:CZ	2.26	0.69
1:A:263:PRO:HB2	1:A:264:PRO:CD	2.22	0.69
1:B:627:ALA:C	1:B:628:ILE:HD12	2.13	0.69
1:C:263:PRO:HB2	1:C:264:PRO:CD	2.22	0.69
1:C:466:LYS:O	1:C:470:ARG:HG3	1.93	0.68
1:A:97:ALA:O	1:A:101:ILE:HG13	1.93	0.68
1:B:622:GLY:O	1:B:626:GLN:HG2	1.93	0.68
1:B:309:SER:HB3	1:B:365:PHE:CZ	2.28	0.68
1:B:263:PRO:HB2	1:B:264:PRO:CD	2.24	0.68
1:C:170:ALA:O	1:C:174:VAL:HG23	1.95	0.67
1:C:238:ARG:HB2	1:C:238:ARG:NH1	2.04	0.67
1:A:622:GLY:O	1:A:626:GLN:HG2	1.95	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:627:ALA:C	1:A:628:ILE:HD12	2.14	0.66
1:B:341:LEU:HD13	1:B:395:ARG:HG2	1.76	0.66
1:A:507:VAL:HG23	1:A:509:ILE:HG12	1.77	0.66
1:C:261:ILE:HG13	1:C:264:PRO:HD2	1.78	0.65
1:C:275:LEU:O	1:C:277:MET:HG3	1.96	0.65
1:A:170:ALA:O	1:A:174:VAL:HG23	1.96	0.65
1:C:345:ILE:HD12	1:C:370:VAL:HG12	1.78	0.65
1:B:369:ASN:HD21	1:B:372:TYR:HB2	1.61	0.64
1:A:254:GLY:HA3	1:A:368:ASP:OD2	1.97	0.64
1:C:507:VAL:HG23	1:C:509:ILE:HG12	1.78	0.64
1:A:531:GLY:O	1:A:577:GLY:HA2	1.98	0.64
1:C:369:ASN:HD21	1:C:372:TYR:HB2	1.61	0.64
1:C:114:LEU:HD11	1:C:197:GLU:HG3	1.77	0.64
1:C:490:GLY:HA3	1:C:495:TYR:CD2	2.33	0.64
1:B:261:ILE:HG13	1:B:264:PRO:HD2	1.79	0.63
1:C:238:ARG:CB	1:C:238:ARG:HH11	2.03	0.63
1:A:106:ASP:O	1:A:111:GLN:HG2	1.99	0.63
1:A:261:ILE:HG13	1:A:264:PRO:HD2	1.80	0.63
1:C:254:GLY:HA3	1:C:368:ASP:OD2	1.97	0.63
1:B:507:VAL:HG23	1:B:509:ILE:HG12	1.81	0.63
1:A:191:GLU:O	1:A:194:ARG:HG2	1.99	0.62
1:B:446:GLU:HB2	4:B:2065:HOH:O	1.98	0.62
1:A:466:LYS:O	1:A:470:ARG:HG3	2.00	0.62
1:A:149:PRO:O	1:A:152:ILE:HG22	2.00	0.62
1:A:545:TYR:HA	1:A:548:LYS:HB2	1.81	0.62
1:B:114:LEU:HD11	1:B:197:GLU:HG3	1.80	0.61
1:C:150:PRO:CG	1:C:182:SER:HB2	2.30	0.61
1:A:113:ALA:O	1:A:117:ILE:HG12	2.00	0.61
1:B:578:VAL:O	1:B:582:VAL:HG23	1.98	0.61
1:A:14:THR:OG1	1:A:582:VAL:HG13	2.00	0.61
1:A:490:GLY:HA3	1:A:495:TYR:CD2	2.35	0.61
1:B:114:LEU:O	1:B:118:GLN:HG3	2.00	0.61
1:C:489:TRP:HB2	4:C:2066:HOH:O	2.01	0.61
1:C:14:THR:OG1	1:C:582:VAL:HG13	2.00	0.60
1:C:263:PRO:HB2	1:C:264:PRO:HD3	1.83	0.60
1:C:191:GLU:O	1:C:194:ARG:HG2	2.01	0.60
1:B:345:ILE:HD12	1:B:370:VAL:HG12	1.83	0.60
1:A:263:PRO:HB2	1:A:264:PRO:HD3	1.81	0.60
1:B:150:PRO:CG	1:B:182:SER:HB2	2.32	0.60
1:A:140:TYR:CD1	1:A:141:PRO:HD2	2.36	0.60
1:B:461:TYR:HD1	1:B:461:TYR:H	1.49	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:369:ASN:HD21	1:A:372:TYR:HB2	1.65	0.60
1:C:108:GLU:HB3	1:C:109:PRO:HD3	1.82	0.60
1:A:150:PRO:CG	1:A:182:SER:HB2	2.32	0.60
1:C:140:TYR:CD1	1:C:141:PRO:HD2	2.36	0.60
1:B:239:ALA:O	1:B:243:ARG:HG2	2.02	0.59
1:A:578:VAL:O	1:A:582:VAL:HG23	2.01	0.59
1:C:270:ILE:O	1:C:274:ILE:HG12	2.02	0.59
1:B:149:PRO:O	1:B:152:ILE:HG22	2.02	0.59
1:B:13:ARG:HH22	1:B:17:ARG:NH2	1.99	0.59
1:B:490:GLY:HA3	1:B:495:TYR:CD2	2.38	0.58
1:C:15:LEU:O	1:C:19:VAL:HG23	2.03	0.58
1:B:140:TYR:CD1	1:B:141:PRO:HD2	2.37	0.58
1:B:263:PRO:HB2	1:B:264:PRO:HD3	1.85	0.58
1:B:43:GLU:OE1	1:B:67:TYR:HE2	1.86	0.58
1:C:92:ILE:HD12	1:C:126:SER:HB3	1.86	0.58
1:B:191:GLU:O	1:B:194:ARG:HG2	2.03	0.58
1:C:149:PRO:O	1:C:152:ILE:HG22	2.04	0.58
1:B:66:ARG:HH21	1:C:115:ARG:HH21	1.52	0.58
1:C:114:LEU:O	1:C:118:GLN:HG3	2.03	0.57
1:C:383:TRP:O	1:C:387:THR:HG23	2.03	0.57
1:A:256:GLY:O	1:A:286:GLY:HA2	2.05	0.57
1:B:92:ILE:HD12	1:B:126:SER:HB3	1.86	0.57
1:A:91:THR:HG22	1:A:117:ILE:HD11	1.84	0.57
1:B:280:HIS:O	1:B:284:ILE:HG12	2.05	0.57
1:C:61:MET:HB3	1:C:101:ILE:HG23	1.85	0.57
1:C:461:TYR:HD1	1:C:461:TYR:H	1.51	0.57
1:A:115:ARG:HH22	1:C:66:ARG:HH21	1.52	0.57
1:A:108:GLU:HB2	1:A:109:PRO:HD3	1.85	0.57
1:A:219:ASP:O	1:A:223:ARG:HG3	2.05	0.56
1:B:466:LYS:O	1:B:470:ARG:HG3	2.05	0.56
1:C:280:HIS:ND1	1:C:281:PRO:HD2	2.20	0.56
1:A:345:ILE:HD12	1:A:370:VAL:HG12	1.85	0.56
1:A:49:LEU:HD11	1:A:53:LEU:HD11	1.88	0.56
1:C:81:TYR:CE2	1:C:537:CYS:HB2	2.40	0.56
1:C:239:ALA:O	1:C:243:ARG:HG2	2.05	0.56
1:C:150:PRO:HG3	1:C:182:SER:HB2	1.87	0.56
1:A:137:VAL:HB	1:A:189:LEU:HG	1.87	0.56
1:B:435:CYS:HB3	1:B:440:VAL:HG21	1.88	0.56
1:C:505:LYS:HD3	1:C:505:LYS:O	2.06	0.56
1:A:150:PRO:HA	1:A:178:SER:HB2	1.87	0.56
1:A:498:GLY:HA2	1:A:562:ALA:HB2	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:150:PRO:HA	1:C:178:SER:HB2	1.88	0.56
1:A:471:ALA:O	1:A:474:TYR:HB3	2.06	0.55
1:B:150:PRO:HA	1:B:178:SER:HB2	1.87	0.55
1:C:392:ASP:OD2	1:C:395:ARG:HB2	2.06	0.55
1:A:150:PRO:HG3	1:A:182:SER:HB2	1.88	0.55
1:B:251:ARG:HH11	1:B:251:ARG:HG3	1.71	0.55
1:C:137:VAL:HB	1:C:189:LEU:HG	1.89	0.55
1:B:137:VAL:HB	1:B:189:LEU:HG	1.89	0.55
1:B:498:GLY:HA2	1:B:562:ALA:HB2	1.89	0.55
1:B:106:ASP:O	1:B:111:GLN:HG2	2.07	0.55
1:B:293:TYR:O	1:B:305:GLN:HG3	2.07	0.55
1:A:505:LYS:HD3	1:A:505:LYS:O	2.06	0.55
1:A:494:LEU:HD22	1:A:559:ALA:HB2	1.87	0.55
1:C:326:LEU:HD22	1:C:330:HIS:CD2	2.41	0.55
1:C:498:GLY:HA2	1:C:562:ALA:HB2	1.88	0.55
1:A:428:LEU:N	1:A:429:PRO:HD2	2.23	0.54
1:A:92:ILE:HD12	1:A:126:SER:HB3	1.89	0.54
1:B:539:SER:HB3	1:B:545:TYR:O	2.08	0.54
1:A:13:ARG:HH11	1:A:13:ARG:HG2	1.73	0.54
1:A:392:ASP:OD2	1:A:395:ARG:HB2	2.07	0.54
1:C:208:ARG:HH11	1:C:208:ARG:HB3	1.73	0.54
1:A:435:CYS:HB3	1:A:440:VAL:HG21	1.89	0.54
1:A:548:LYS:HD2	4:A:2072:HOH:O	2.08	0.54
1:B:238:ARG:HB2	1:B:238:ARG:HH11	1.73	0.54
1:B:81:TYR:CE2	1:B:537:CYS:HB2	2.43	0.54
1:C:471:ALA:O	1:C:474:TYR:HB3	2.08	0.53
1:C:494:LEU:HD22	1:C:559:ALA:HB2	1.89	0.53
1:B:574:ALA:O	1:B:578:VAL:HG23	2.09	0.53
1:C:14:THR:HG23	1:C:585:GLN:OE1	2.08	0.53
1:A:147:MET:HG2	1:A:225:LEU:HB3	1.90	0.53
1:A:238:ARG:CB	1:A:238:ARG:HH11	2.16	0.53
1:A:293:TYR:O	1:A:305:GLN:HG3	2.08	0.53
1:B:536:ASP:OD2	1:B:538:ARG:HG3	2.09	0.53
1:C:297:LEU:HB2	1:C:301:GLY:O	2.09	0.53
1:A:312:TRP:CZ3	1:A:380:VAL:HG21	2.44	0.53
1:A:536:ASP:OD2	1:A:538:ARG:HG3	2.09	0.53
1:B:15:LEU:O	1:B:19:VAL:HG23	2.09	0.53
1:C:542:ASP:OD2	1:C:544:ALA:HB3	2.08	0.53
1:A:61:MET:HB3	1:A:101:ILE:HG23	1.90	0.53
1:C:578:VAL:O	1:C:582:VAL:HG23	2.07	0.53
1:A:81:TYR:CE2	1:A:537:CYS:HB2	2.44	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:MET:HB2	1:A:226:HIS:CE1	2.44	0.53
1:A:403:GLY:O	1:A:407:ILE:HG13	2.09	0.53
1:B:150:PRO:HG3	1:B:182:SER:HB2	1.90	0.52
1:C:428:LEU:N	1:C:429:PRO:HD2	2.24	0.52
1:B:208:ARG:HB3	1:B:208:ARG:HH11	1.73	0.52
1:C:345:ILE:HD11	4:C:2053:HOH:O	2.09	0.52
1:A:81:TYR:HB2	1:A:82:PRO:HD2	1.92	0.52
1:A:300:GLY:HA3	4:A:2011:HOH:O	2.10	0.52
1:A:383:TRP:CH2	1:A:454:GLU:HG2	2.45	0.52
1:B:576:ARG:HG3	1:B:576:ARG:HH11	1.75	0.52
1:B:61:MET:HB3	1:B:101:ILE:HG23	1.92	0.52
1:B:73:ARG:HD3	1:B:77:THR:O	2.10	0.52
1:C:312:TRP:CZ3	1:C:380:VAL:HG21	2.44	0.52
1:B:312:TRP:CZ3	1:B:380:VAL:HG21	2.45	0.52
1:C:185:PRO:HG3	1:C:274:ILE:CD1	2.40	0.52
1:C:435:CYS:HB3	1:C:440:VAL:HG21	1.91	0.51
1:C:81:TYR:HB2	1:C:82:PRO:HD2	1.92	0.51
1:B:81:TYR:HB2	1:B:82:PRO:HD2	1.93	0.51
1:A:208:ARG:HB3	1:A:208:ARG:HH11	1.75	0.51
1:A:278:THR:HA	1:A:283:PHE:CD2	2.45	0.51
1:A:473:GLU:O	1:A:477:ARG:HG3	2.09	0.51
1:B:394:ARG:HA	1:B:394:ARG:NH1	2.26	0.51
1:B:217:ILE:H	1:B:217:ILE:CD1	2.14	0.51
1:B:376:ASP:HB3	3:B:800:R23:H4E1	1.91	0.51
1:C:13:ARG:HH22	1:C:17:ARG:NH2	2.09	0.51
4:B:2002:HOH:O	1:C:548:LYS:HA	2.10	0.51
1:B:147:MET:HG2	1:B:225:LEU:HB3	1.93	0.51
3:B:800:R23:H3E1	4:B:2084:HOH:O	2.10	0.51
1:C:574:ALA:O	1:C:578:VAL:HG23	2.10	0.51
1:B:254:GLY:HA3	1:B:368:ASP:OD2	2.10	0.51
1:A:270:ILE:O	1:A:274:ILE:HG12	2.12	0.50
1:A:574:ALA:O	1:A:578:VAL:HG23	2.11	0.50
1:B:219:ASP:O	1:B:223:ARG:HG3	2.10	0.50
1:A:158:ARG:O	1:A:158:ARG:HG2	2.11	0.50
1:C:383:TRP:CH2	1:C:454:GLU:HG2	2.47	0.50
1:C:322:ARG:NH2	1:C:328:ALA:HB2	2.27	0.50
1:C:536:ASP:OD2	1:C:538:ARG:HG3	2.12	0.50
1:B:283:PHE:CD1	1:B:283:PHE:C	2.85	0.50
1:A:238:ARG:NH1	1:A:238:ARG:HB2	2.17	0.50
1:B:471:ALA:O	1:B:474:TYR:HB3	2.12	0.50
1:C:52:ILE:HD12	1:C:187:PHE:CD2	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:560:LEU:O	1:A:564:ILE:HG13	2.12	0.50
1:B:52:ILE:HD12	1:B:187:PHE:CD2	2.46	0.50
1:B:322:ARG:NH2	1:B:328:ALA:HB2	2.27	0.50
1:A:128:VAL:HG13	1:A:146:PRO:HD2	1.94	0.49
1:A:576:ARG:HH11	1:A:576:ARG:HG3	1.77	0.49
1:C:345:ILE:HD12	1:C:370:VAL:HA	1.94	0.49
1:B:428:LEU:N	1:B:429:PRO:HD2	2.27	0.49
1:B:473:GLU:O	1:B:477:ARG:HG3	2.12	0.49
1:C:411:GLN:HB2	1:C:417:TRP:CZ2	2.46	0.49
1:C:280:HIS:O	1:C:284:ILE:HG13	2.13	0.49
1:C:158:ARG:HG2	1:C:158:ARG:O	2.12	0.49
1:C:147:MET:HG2	1:C:225:LEU:HB3	1.95	0.49
1:A:345:ILE:HD12	1:A:370:VAL:HA	1.93	0.49
1:A:378:THR:O	1:A:382:VAL:HG23	2.13	0.49
1:C:128:VAL:HG13	1:C:146:PRO:HD2	1.95	0.49
1:A:14:THR:HG23	1:A:585:GLN:CD	2.33	0.49
1:B:52:ILE:HG23	1:B:187:PHE:CD2	2.48	0.49
1:C:66:ARG:HG2	1:C:66:ARG:HH11	1.77	0.49
1:A:554:SER:OG	1:A:597:THR:HG21	2.12	0.49
1:C:544:ALA:O	1:C:548:LYS:HD3	2.12	0.49
1:B:128:VAL:HG13	1:B:146:PRO:HD2	1.95	0.49
1:B:152:ILE:O	1:B:152:ILE:HG12	2.13	0.49
1:C:576:ARG:HG3	1:C:576:ARG:HH11	1.78	0.49
1:C:65:ARG:O	1:C:69:LEU:HD23	2.13	0.49
1:C:45:GLU:HG2	1:C:133:TRP:CE2	2.48	0.48
1:C:86:PRO:HB2	1:C:116:PHE:CZ	2.48	0.48
1:A:513:GLU:HB2	1:A:516:ILE:HG12	1.95	0.48
1:B:158:ARG:O	1:B:158:ARG:HG2	2.12	0.48
1:B:49:LEU:HD11	1:B:53:LEU:HD11	1.95	0.48
1:B:554:SER:OG	1:B:597:THR:HG21	2.13	0.48
1:B:392:ASP:OD2	1:B:395:ARG:HB2	2.13	0.48
1:A:52:ILE:HD12	1:A:187:PHE:CD2	2.47	0.48
1:B:270:ILE:O	1:B:274:ILE:HG12	2.13	0.48
1:B:108:GLU:HB3	1:B:109:PRO:HD3	1.96	0.48
1:B:77:THR:OG1	1:B:86:PRO:HB3	2.13	0.48
1:C:404:PHE:C	1:C:404:PHE:CD1	2.87	0.48
1:B:370:VAL:N	4:B:2048:HOH:O	2.47	0.48
1:B:383:TRP:CH2	1:B:454:GLU:HG2	2.49	0.48
1:C:473:GLU:O	1:C:477:ARG:HG3	2.14	0.48
1:A:52:ILE:HG23	1:A:187:PHE:CD2	2.49	0.48
1:C:560:LEU:O	1:C:564:ILE:HG13	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:GLU:HG2	1:A:133:TRP:CE2	2.49	0.47
1:B:505:LYS:O	1:B:505:LYS:HD3	2.14	0.47
1:B:580:TYR:O	1:B:584:THR:HG23	2.14	0.47
1:C:613:ARG:O	1:C:617:PRO:HG2	2.13	0.47
1:A:255:ASP:HB3	4:A:2038:HOH:O	2.14	0.47
1:C:152:ILE:HG12	1:C:152:ILE:O	2.14	0.47
1:B:245:LEU:HD22	1:B:275:LEU:HD13	1.97	0.47
1:B:345:ILE:HD12	1:B:370:VAL:HA	1.95	0.47
1:B:494:LEU:HD22	1:B:559:ALA:HB2	1.95	0.47
1:C:280:HIS:CE1	1:C:281:PRO:HD2	2.49	0.47
1:C:461:TYR:N	1:C:461:TYR:CD1	2.82	0.47
1:A:262:GLN:HB3	1:A:263:PRO:HD3	1.96	0.47
1:A:542:ASP:OD2	1:A:544:ALA:HB3	2.15	0.47
1:B:13:ARG:HH11	1:B:13:ARG:HG2	1.80	0.47
1:C:332:ARG:HG2	1:C:332:ARG:HH11	1.80	0.47
1:A:613:ARG:O	1:A:617:PRO:HG2	2.15	0.47
1:B:251:ARG:NH1	1:B:251:ARG:HG3	2.29	0.47
1:C:185:PRO:HG3	1:C:274:ILE:HD11	1.97	0.47
1:B:143:GLU:N	1:B:143:GLU:OE1	2.44	0.47
1:B:14:THR:OG1	1:B:582:VAL:HG13	2.15	0.47
1:B:29:GLU:H	1:B:29:GLU:CD	2.18	0.47
1:B:192:ARG:HH11	1:B:192:ARG:HG2	1.80	0.47
1:C:43:GLU:OE1	1:C:67:TYR:HE2	1.98	0.46
1:C:143:GLU:N	1:C:143:GLU:OE1	2.47	0.46
1:A:322:ARG:NH2	1:A:328:ALA:HB2	2.31	0.46
1:C:52:ILE:HG23	1:C:187:PHE:CD2	2.50	0.46
1:C:531:GLY:O	1:C:577:GLY:HA2	2.16	0.46
1:A:11:TYR:CD1	1:A:12:ALA:N	2.83	0.46
1:B:31:TYR:HB2	1:B:314:THR:OG1	2.16	0.46
1:A:405:ARG:HH11	1:A:405:ARG:HG2	1.80	0.46
1:A:92:ILE:O	1:A:96:VAL:HG23	2.16	0.46
1:C:586:ARG:HD3	1:C:592:ASP:OD2	2.15	0.46
1:A:81:TYR:CD1	1:A:81:TYR:N	2.84	0.46
1:B:326:LEU:HD22	1:B:330:HIS:CD2	2.51	0.45
1:B:461:TYR:N	1:B:461:TYR:CD1	2.84	0.45
1:C:219:ASP:O	1:C:223:ARG:HG3	2.16	0.45
1:C:29:GLU:H	1:C:29:GLU:CD	2.19	0.45
1:C:403:GLY:O	1:C:407:ILE:HG12	2.16	0.45
1:C:513:GLU:HB2	1:C:516:ILE:HG12	1.97	0.45
1:C:106:ASP:O	1:C:111:GLN:HG2	2.16	0.45
1:C:31:TYR:HB2	1:C:314:THR:OG1	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:49:LEU:HD11	1:C:53:LEU:HD11	1.98	0.45
1:C:113:ALA:O	1:C:117:ILE:HG13	2.16	0.45
1:C:279:GLN:HE21	1:C:279:GLN:HA	1.81	0.45
1:A:332:ARG:HG2	1:A:332:ARG:HH11	1.82	0.45
1:C:293:TYR:O	1:C:305:GLN:HG3	2.15	0.45
1:A:261:ILE:HG13	1:A:264:PRO:CD	2.45	0.45
1:B:261:ILE:HG13	1:B:264:PRO:CD	2.46	0.45
1:B:53:LEU:O	1:B:55:ARG:HG3	2.17	0.45
1:C:192:ARG:HH11	1:C:192:ARG:HG2	1.82	0.45
1:C:554:SER:OG	1:C:597:THR:HG21	2.16	0.45
1:A:143:GLU:N	1:A:143:GLU:OE1	2.45	0.45
1:C:114:LEU:O	1:C:114:LEU:HD23	2.16	0.45
1:C:378:THR:O	1:C:382:VAL:HG23	2.17	0.45
1:C:99:LYS:HD3	1:C:195:VAL:HG12	1.99	0.45
1:B:43:GLU:OE1	1:B:67:TYR:CE2	2.70	0.45
1:C:13:ARG:HG2	1:C:13:ARG:HH11	1.82	0.45
1:A:152:ILE:HG12	1:A:152:ILE:O	2.16	0.44
1:A:29:GLU:H	1:A:29:GLU:CD	2.19	0.44
1:B:332:ARG:HH11	1:B:332:ARG:HG2	1.81	0.44
1:A:228:TYR:HD1	1:A:231:LEU:HD12	1.82	0.44
1:C:295:VAL:HB	1:C:303:MET:HG2	1.97	0.44
1:A:31:TYR:HB2	1:A:314:THR:OG1	2.17	0.44
1:B:208:ARG:NH1	1:B:208:ARG:HB3	2.32	0.44
1:B:262:GLN:HB3	1:B:263:PRO:HD3	1.99	0.44
1:B:542:ASP:OD2	1:B:544:ALA:HB3	2.17	0.44
1:C:261:ILE:HG13	1:C:264:PRO:CD	2.46	0.44
1:C:555:GLN:HA	1:C:558:TRP:CE3	2.52	0.44
1:A:86:PRO:HB2	1:A:116:PHE:CE2	2.51	0.44
1:A:251:ARG:HG3	1:A:251:ARG:HH11	1.82	0.44
1:A:510:ASP:OD2	1:A:512:ARG:HB3	2.17	0.44
1:B:555:GLN:HA	1:B:558:TRP:CE3	2.53	0.44
1:C:405:ARG:HH11	1:C:405:ARG:HG2	1.83	0.44
1:A:310:PRO:HD2	1:A:366:GLN:HA	2.00	0.44
1:C:150:PRO:HG2	1:C:182:SER:HB2	1.99	0.44
1:C:208:ARG:HB3	1:C:208:ARG:NH1	2.31	0.44
1:C:509:ILE:HG22	1:C:510:ASP:N	2.32	0.44
1:C:92:ILE:O	1:C:96:VAL:HG23	2.18	0.44
1:A:192:ARG:HH11	1:A:192:ARG:HG2	1.82	0.44
1:A:221:LEU:O	1:A:221:LEU:HD12	2.17	0.44
1:A:555:GLN:HA	1:A:558:TRP:CE3	2.53	0.44
1:A:613:ARG:HG3	1:A:614:HIS:CE1	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:509:ILE:HG22	1:B:510:ASP:N	2.32	0.44
1:C:334:VAL:HG13	1:C:390:LEU:HD22	2.00	0.44
1:A:199:TYR:HA	4:A:2032:HOH:O	2.17	0.44
1:B:45:GLU:HG2	1:B:133:TRP:CE2	2.53	0.44
1:B:418:GLY:O	1:B:448:VAL:HG11	2.18	0.44
1:B:221:LEU:O	1:B:224:ALA:HB3	2.18	0.44
1:B:366:GLN:CD	1:B:366:GLN:H	2.20	0.44
1:C:418:GLY:O	1:C:448:VAL:HG11	2.17	0.44
1:A:49:LEU:HD22	1:A:266:PHE:CD2	2.53	0.43
1:A:533:TRP:HB2	1:A:556:THR:HA	2.00	0.43
1:B:37:LEU:N	1:B:37:LEU:HD22	2.33	0.43
1:C:283:PHE:C	1:C:283:PHE:CD1	2.91	0.43
1:C:43:GLU:OE1	1:C:67:TYR:CE2	2.70	0.43
1:B:108:GLU:HG2	1:B:112:LYS:HE3	2.00	0.43
1:B:92:ILE:O	1:B:96:VAL:HG23	2.18	0.43
1:C:535:GLU:HA	1:C:548:LYS:O	2.18	0.43
1:B:533:TRP:HB2	1:B:556:THR:HA	2.00	0.43
1:A:295:VAL:HB	1:A:303:MET:HG2	1.99	0.43
1:A:406:TRP:O	1:A:410:MET:HG2	2.18	0.43
1:B:11:TYR:CD1	1:B:12:ALA:N	2.87	0.43
1:B:150:PRO:HG2	1:B:182:SER:HB2	2.00	0.43
1:B:192:ARG:HG2	1:B:192:ARG:NH1	2.33	0.43
1:C:357:ASN:N	1:C:357:ASN:HD22	2.15	0.43
1:A:348:PRO:HG2	2:A:700:C8E:H71	2.01	0.43
1:B:560:LEU:O	1:B:564:ILE:HG13	2.18	0.43
1:C:366:GLN:CD	1:C:366:GLN:H	2.20	0.43
1:C:613:ARG:HG3	1:C:614:HIS:CE1	2.53	0.43
1:C:322:ARG:HH21	1:C:328:ALA:HB2	1.83	0.43
1:C:326:LEU:HA	1:C:327:PRO:HD3	1.87	0.43
1:C:404:PHE:O	1:C:408:VAL:HG23	2.18	0.43
1:B:357:ASN:N	1:B:357:ASN:HD22	2.15	0.43
1:B:520:LEU:HD12	1:B:520:LEU:N	2.34	0.43
1:C:43:GLU:CD	1:C:43:GLU:H	2.22	0.43
1:A:370:VAL:N	4:A:2050:HOH:O	2.52	0.43
1:A:377:ASP:O	1:A:381:VAL:HG23	2.18	0.43
1:A:509:ILE:HG22	1:A:510:ASP:N	2.33	0.43
1:A:418:GLY:O	1:A:448:VAL:HG11	2.19	0.43
1:B:613:ARG:O	1:B:617:PRO:HG2	2.18	0.43
1:B:305:GLN:HA	4:B:2006:HOH:O	2.19	0.42
1:B:378:THR:O	1:B:382:VAL:HG23	2.20	0.42
1:B:81:TYR:N	1:B:81:TYR:CD1	2.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:11:TYR:CD1	1:C:12:ALA:N	2.88	0.42
1:A:357:ASN:N	1:A:357:ASN:HD22	2.16	0.42
1:C:53:LEU:O	1:C:55:ARG:HG3	2.18	0.42
1:A:154:PHE:CE2	1:A:240:ALA:HB2	2.54	0.42
1:A:86:PRO:HB2	1:A:116:PHE:CZ	2.53	0.42
1:B:529:PRO:C	1:B:531:GLY:H	2.23	0.42
1:B:576:ARG:NH1	1:B:576:ARG:HG3	2.34	0.42
1:C:251:ARG:HH11	1:C:251:ARG:HG3	1.84	0.42
1:A:394:ARG:O	1:A:394:ARG:HD3	2.19	0.42
1:A:42:MET:SD	1:A:262:GLN:HG3	2.59	0.42
1:B:191:GLU:HA	1:B:194:ARG:HG2	2.01	0.42
1:B:256:GLY:O	1:B:286:GLY:HA2	2.19	0.42
1:B:513:GLU:HB2	1:B:516:ILE:HG12	2.01	0.42
1:B:243:ARG:HD2	4:B:2029:HOH:O	2.18	0.42
1:C:394:ARG:NH1	1:C:394:ARG:HA	2.35	0.42
1:A:192:ARG:NH1	1:A:192:ARG:HG2	2.35	0.42
1:A:40:VAL:HG13	4:A:2006:HOH:O	2.20	0.42
1:B:404:PHE:C	1:B:404:PHE:CD1	2.93	0.42
1:B:557:ALA:O	1:B:561:MET:HG3	2.19	0.42
1:C:104:SER:C	1:C:106:ASP:H	2.22	0.42
1:C:81:TYR:CD1	1:C:81:TYR:N	2.87	0.42
1:A:150:PRO:HG2	1:A:182:SER:HB2	2.01	0.42
1:A:208:ARG:HB3	1:A:208:ARG:NH1	2.33	0.42
1:B:405:ARG:HG2	1:B:405:ARG:HH11	1.85	0.42
1:B:48:LEU:O	1:B:52:ILE:HG12	2.20	0.42
1:C:37:LEU:HD22	1:C:37:LEU:N	2.34	0.42
1:C:628:ILE:HD12	1:C:628:ILE:N	2.35	0.42
1:A:322:ARG:HD3	1:A:333:LEU:HD13	2.01	0.42
1:C:147:MET:HB2	1:C:226:HIS:CE1	2.55	0.42
1:A:347:VAL:HA	1:A:348:PRO:HD3	1.91	0.41
1:A:617:PRO:O	1:A:621:LEU:HG	2.20	0.41
1:A:394:ARG:O	1:A:398:ASP:HB2	2.20	0.41
1:B:332:ARG:HD2	1:B:332:ARG:HA	1.88	0.41
1:B:279:GLN:HA	1:B:284:ILE:HD11	2.01	0.41
1:C:186:VAL:O	1:C:188:PRO:HD3	2.20	0.41
1:C:262:GLN:HB3	1:C:263:PRO:HD3	2.01	0.41
1:C:617:PRO:O	1:C:621:LEU:HG	2.20	0.41
1:A:191:GLU:HA	1:A:194:ARG:HG2	2.01	0.41
1:A:429:PRO:O	1:A:432:ILE:HG12	2.21	0.41
1:A:448:VAL:O	1:A:452:VAL:HG23	2.20	0.41
1:B:42:MET:SD	1:B:262:GLN:HG3	2.60	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:295:VAL:HB	1:B:303:MET:HG2	2.01	0.41
1:C:310:PRO:HD2	1:C:366:GLN:HA	2.01	0.41
1:A:405:ARG:NH1	1:A:405:ARG:HG2	2.35	0.41
1:A:492:ASN:HB3	1:A:535:GLU:CD	2.41	0.41
1:A:14:THR:HG23	1:A:585:GLN:OE1	2.20	0.41
1:B:347:VAL:HA	1:B:348:PRO:HD3	1.92	0.41
1:B:386:ASN:HA	1:B:396:ARG:HH12	1.85	0.41
1:C:490:GLY:HA3	1:C:495:TYR:CE2	2.55	0.41
1:C:533:TRP:HB2	1:C:556:THR:HA	2.02	0.41
1:B:217:ILE:N	1:B:217:ILE:HD12	2.20	0.41
1:B:392:ASP:HB3	4:B:2052:HOH:O	2.20	0.41
1:B:490:GLY:HA3	1:B:495:TYR:CE2	2.56	0.41
1:C:192:ARG:NH1	1:C:192:ARG:HG2	2.35	0.41
1:C:359:LYS:HA	1:C:360:PRO:HD3	1.90	0.41
1:A:108:GLU:OE2	1:C:66:ARG:NH1	2.53	0.41
1:A:153:MET:HG3	1:A:244:ALA:HB2	2.03	0.41
1:A:245:LEU:HD22	1:A:275:LEU:HD13	2.02	0.41
1:A:92:ILE:HG12	1:A:117:ILE:HD12	2.03	0.41
1:C:191:GLU:HA	1:C:194:ARG:HG2	2.02	0.41
1:B:124:GLU:OE2	1:B:140:TYR:HE1	2.03	0.41
1:C:405:ARG:HG2	1:C:405:ARG:NH1	2.35	0.41
1:B:322:ARG:HH21	1:B:328:ALA:HB2	1.86	0.41
1:C:510:ASP:OD2	1:C:512:ARG:HB3	2.20	0.41
1:A:114:LEU:HD11	1:A:197:GLU:HG3	2.03	0.41
1:A:88:LEU:HG	1:A:92:ILE:HD11	2.03	0.41
1:A:292:LEU:HD12	1:A:292:LEU:O	2.21	0.41
1:A:396:ARG:O	1:A:400:MET:HG3	2.19	0.41
1:B:99:LYS:HD3	1:B:195:VAL:HG12	2.02	0.41
1:C:118:GLN:NE2	1:C:200:GLU:O	2.54	0.41
1:A:359:LYS:HA	1:A:360:PRO:HD3	1.93	0.40
1:A:43:GLU:CD	1:A:43:GLU:H	2.24	0.40
1:A:490:GLY:HA3	1:A:495:TYR:CE2	2.56	0.40
1:A:310:PRO:HG2	1:A:311:VAL:H	1.86	0.40
1:C:275:LEU:O	1:C:276:ASP:C	2.60	0.40
1:A:520:LEU:N	1:A:520:LEU:HD12	2.37	0.40
1:A:529:PRO:C	1:A:531:GLY:H	2.25	0.40
1:A:576:ARG:NH1	1:A:576:ARG:HG3	2.36	0.40
1:B:149:PRO:HA	1:B:150:PRO:HD3	1.84	0.40
1:B:617:PRO:O	1:B:621:LEU:HG	2.21	0.40
1:C:11:TYR:O	1:C:14:THR:HB	2.21	0.40
1:C:310:PRO:HG2	1:C:311:VAL:H	1.85	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:610:THR:HG22	1:C:611:MET:CE	2.50	0.40
1:B:186:VAL:O	1:B:188:PRO:HD3	2.22	0.40
1:C:208:ARG:CB	1:C:208:ARG:NH1	2.85	0.40
1:A:612:TYR:HB2	4:A:2077:HOH:O	2.22	0.40
1:B:520:LEU:HD12	1:B:520:LEU:H	1.86	0.40
1:C:280:HIS:ND1	1:C:281:PRO:CD	2.84	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	618/631 (98%)	550 (89%)	62 (10%)	6 (1%)	18	50
1	B	618/631 (98%)	553 (90%)	58 (9%)	7 (1%)	17	47
1	C	618/631 (98%)	555 (90%)	56 (9%)	7 (1%)	17	47
All	All	1854/1893 (98%)	1658 (89%)	176 (10%)	20 (1%)	17	47

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	107	GLU
1	C	105	ARG
1	A	200	GLU
1	A	263	PRO
1	A	361	GLY
1	B	263	PRO
1	B	361	GLY
1	C	263	PRO
1	B	200	GLU
1	B	216	TRP
1	C	200	GLU

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Mol	Chain	Res	Type
1	C	216	TRP
1	C	361	GLY
1	A	492	ASN
1	B	492	ASN
1	C	492	ASN
1	B	509	ILE
1	A	509	ILE
1	C	509	ILE
1	B	310	PRO

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	503/513 (98%)	491 (98%)	12 (2%)	54	85
1	B	503/513 (98%)	490 (97%)	13 (3%)	51	83
1	C	503/513 (98%)	490 (97%)	13 (3%)	51	83
All	All	1509/1539 (98%)	1471 (98%)	38 (2%)	53	84

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	LEU
1	A	75	ASP
1	A	115	ARG
1	A	158	ARG
1	A	238	ARG
1	A	258	TRP
1	A	366	GLN
1	A	428	LEU
1	A	447	ASP
1	A	485	TRP
1	A	542	ASP
1	A	591	TRP
1	B	58	ARG

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Mol	Chain	Res	Type
1	B	73	ARG
1	B	115	ARG
1	B	158	ARG
1	B	238	ARG
1	B	366	GLN
1	B	394	ARG
1	B	404	PHE
1	B	428	LEU
1	B	447	ASP
1	B	485	TRP
1	B	542	ASP
1	B	591	TRP
1	C	58	ARG
1	C	75	ASP
1	C	115	ARG
1	C	158	ARG
1	C	238	ARG
1	C	279	GLN
1	C	366	GLN
1	C	394	ARG
1	C	428	LEU
1	C	447	ASP
1	C	485	TRP
1	C	542	ASP
1	C	591	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	357	ASN
1	A	579	GLN
1	B	70	HIS
1	B	279	GLN
1	B	357	ASN
1	B	579	GLN
1	C	279	GLN
1	C	357	ASN
1	C	369	ASN
1	C	386	ASN
1	C	579	GLN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

6 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	C8E	A	700	-	20,20,20	0.85	0	19,19,19	1.76	5 (26%)
3	R23	A	800	-	34,34,34	1.80	11 (32%)	44,46,46	2.07	5 (11%)
2	C8E	B	700	-	20,20,20	0.80	0	19,19,19	1.78	6 (31%)
3	R23	B	800	-	34,34,34	1.76	11 (32%)	44,46,46	2.16	4 (9%)
2	C8E	C	700	-	20,20,20	0.83	0	19,19,19	1.78	6 (31%)
3	R23	C	800	-	34,34,34	1.80	12 (35%)	44,46,46	2.09	5 (11%)

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	C8E	A	700	-	-	0/18/18/18	0/0/0/0
3	R23	A	800	-	-	0/20/38/38	0/4/4/4
2	C8E	B	700	-	-	0/18/18/18	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	R23	B	800	-	-	0/20/38/38	0/4/4/4
2	C8E	C	700	-	-	0/18/18/18	0/0/0/0
3	R23	C	800	-	-	0/20/38/38	0/4/4/4

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	800	R23	C6F-C1F	2.05	1.42	1.38
3	C	800	R23	C2E-N1E	2.06	1.52	1.47
3	A	800	R23	C6F-C1F	2.11	1.43	1.38
3	B	800	R23	C6F-C1F	2.20	1.43	1.38
3	C	800	R23	C2F-C1F	2.20	1.43	1.38
3	B	800	R23	C2C-C1C	2.21	1.42	1.39
3	C	800	R23	C5F-C6F	2.23	1.42	1.38
3	C	800	R23	C2F-C3F	2.24	1.42	1.38
3	A	800	R23	C3C-C4C	2.25	1.43	1.38
3	B	800	R23	C3C-C4C	2.27	1.43	1.38
3	A	800	R23	C6C-C1C	2.29	1.42	1.39
3	C	800	R23	C5F-C4F	2.30	1.42	1.38
3	A	800	R23	C2F-C1F	2.32	1.43	1.38
3	B	800	R23	C2F-C1F	2.33	1.43	1.38
3	A	800	R23	C5F-C6F	2.36	1.43	1.38
3	A	800	R23	C5F-C4F	2.37	1.42	1.38
3	B	800	R23	C6C-C1C	2.39	1.42	1.39
3	C	800	R23	C3C-C4C	2.42	1.43	1.38
3	C	800	R23	C6C-C1C	2.42	1.43	1.39
3	B	800	R23	C2F-C3F	2.50	1.43	1.38
3	B	800	R23	C2C-C3C	2.53	1.43	1.38
3	B	800	R23	C5F-C4F	2.53	1.42	1.38
3	A	800	R23	C2F-C3F	2.54	1.43	1.38
3	B	800	R23	C5F-C6F	2.56	1.43	1.38
3	C	800	R23	C3F-C4F	2.61	1.43	1.38
3	A	800	R23	C3F-C4F	2.68	1.43	1.38
3	A	800	R23	C2C-C3C	2.73	1.43	1.38
3	C	800	R23	C2C-C1C	2.73	1.43	1.39
3	A	800	R23	C2C-C1C	2.79	1.43	1.39
3	B	800	R23	C3F-C4F	2.80	1.43	1.38
3	C	800	R23	C2C-C3C	2.88	1.43	1.38
3	A	800	R23	C1-N4	3.03	1.41	1.36
3	C	800	R23	C1-N4	3.12	1.41	1.36
3	B	800	R23	C1-N4	3.30	1.41	1.36

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	800	R23	O3-C1-O2	-5.34	113.22	123.68
3	B	800	R23	O3-C1-O2	-5.31	113.27	123.68
3	A	800	R23	O3-C1-O2	-5.19	113.52	123.68
2	B	700	C8E	C19-O18-C17	-2.69	101.65	113.30
2	C	700	C8E	C19-O18-C17	-2.65	101.84	113.30
2	A	700	C8E	C19-O18-C17	-2.64	101.86	113.30
2	B	700	C8E	C16-O15-C14	-2.16	103.95	113.30
2	C	700	C8E	C7-C6-C5	-2.13	103.48	114.45
2	C	700	C8E	C16-O15-C14	-2.11	104.18	113.30
2	A	700	C8E	C16-O15-C14	-2.08	104.29	113.30
2	B	700	C8E	C7-C6-C5	-2.07	103.80	114.45
3	A	800	R23	C1A-N4-C1B	2.02	120.13	117.84
3	C	800	R23	C1A-N4-C1B	2.39	120.55	117.84
3	A	800	R23	C2E-N1E-C6E	2.43	114.37	108.87
3	C	800	R23	C2E-N1E-C6E	2.55	114.66	108.87
3	B	800	R23	C2E-N1E-C6E	2.57	114.70	108.87
2	C	700	C8E	O18-C19-C20	2.58	122.07	110.15
2	B	700	C8E	O18-C19-C20	2.59	122.10	110.15
2	A	700	C8E	O18-C19-C20	2.64	122.33	110.15
2	A	700	C8E	O15-C16-C17	2.87	123.60	110.41
2	C	700	C8E	O15-C16-C17	2.89	123.69	110.41
2	B	700	C8E	O15-C16-C17	2.90	123.71	110.41
3	A	800	R23	C1F-O3-C1	3.13	123.01	117.22
3	B	800	R23	C1F-O3-C1	3.33	123.37	117.22
3	C	800	R23	C1F-O3-C1	3.33	123.38	117.22
2	A	700	C8E	O15-C14-C13	4.53	131.20	110.41
2	C	700	C8E	O15-C14-C13	4.55	131.28	110.41
2	B	700	C8E	O15-C14-C13	4.55	131.31	110.41
3	C	800	R23	O3-C1-N4	10.98	123.59	111.88
3	A	800	R23	O3-C1-N4	11.18	123.80	111.88
3	B	800	R23	O3-C1-N4	11.61	124.26	111.88

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	700	C8E	1	0
3	B	800	R23	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 [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, 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	620/631 (98%)	-0.66	0 100 100	6, 29, 81, 99	0
1	B	620/631 (98%)	-0.73	0 100 100	6, 18, 67, 99	0
1	C	620/631 (98%)	-0.72	0 100 100	6, 21, 71, 99	0
All	All	1860/1893 (98%)	-0.70	0 100 100	6, 22, 75, 99	0

There are no RSRZ outliers to report.

### 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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	C8E	C	700	21/21	0.81	0.40	7.50	76,76,76,76	0
2	C8E	B	700	21/21	0.85	0.43	7.47	81,81,81,81	0
2	C8E	A	700	21/21	0.78	0.45	6.35	83,83,83,83	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	R23	A	800	31/31	0.95	0.19	2.42	6,15,26,26	0
3	R23	C	800	31/31	0.94	0.17	1.90	8,13,19,19	0
3	R23	B	800	31/31	0.96	0.17	1.86	6,6,6,6	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.