



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 01:19 pm GMT

PDB ID : 2Z55  
Title : Bacterioruberin in the trimeric structure of archaerhodopsin-2  
Authors : Kouyama, T.; Yoshimura, K.  
Deposited on : 2007-06-28  
Resolution : 2.50 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



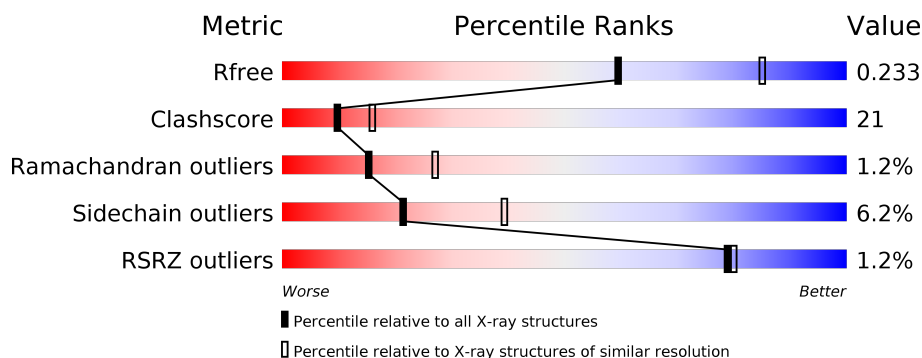
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	3846 (2.50-2.50)
Clashscore	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)
RSRZ outliers	101464	3876 (2.50-2.50)

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	253	
1	B	253	
1	D	253	
1	E	253	

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 crite-



ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MAN	A	282	-	-	-	X
2	GAL	A	283	-	-	-	X
2	MAN	B	282	-	-	-	X
2	GAL	B	283	-	-	-	X
2	MAN	D	282	-	-	-	X
2	GAL	D	283	-	-	-	X
3	RET	D	260	-	-	-	X
3	RET	E	260	-	-	-	X
4	22B	A	270	-	-	-	X
4	22B	B	270	-	-	-	X
4	22B	D	270	-	-	-	X
4	22B	E	270	-	-	-	X
5	L2P	A	280	-	-	-	X
5	L2P	B	280	-	-	-	X
5	L2P	D	280	-	-	-	X
5	L2P	E	280	-	-	-	X



## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8087 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 Archaelhodopsin-2.

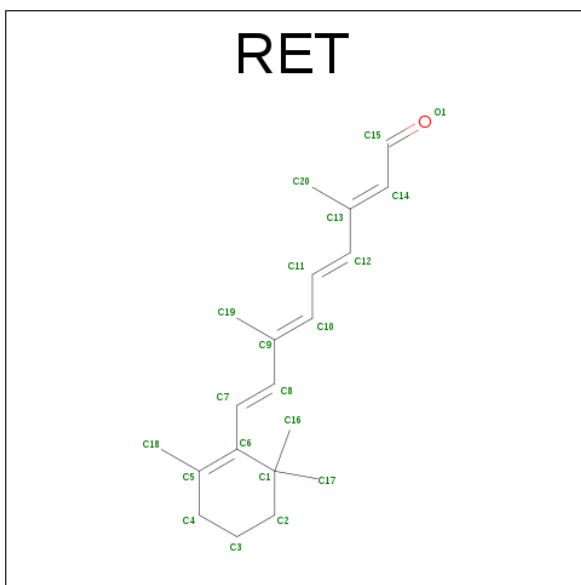
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	242	Total	C	N	O	S	0	0	0
			1868	1243	290	331	4			
1	B	238	Total	C	N	O	S	0	0	0
			1838	1225	286	323	4			
1	D	237	Total	C	N	O	S	0	0	0
			1830	1219	284	323	4			
1	E	238	Total	C	N	O	S	0	0	0
			1838	1225	286	323	4			

- Molecule 2 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	3	Total	C	O	0	0
			33	18	15		
2	B	3	Total	C	O	0	0
			33	18	15		
2	D	3	Total	C	O	0	0
			33	18	15		
2	D	3	Total	C	O	0	0
			33	18	15		

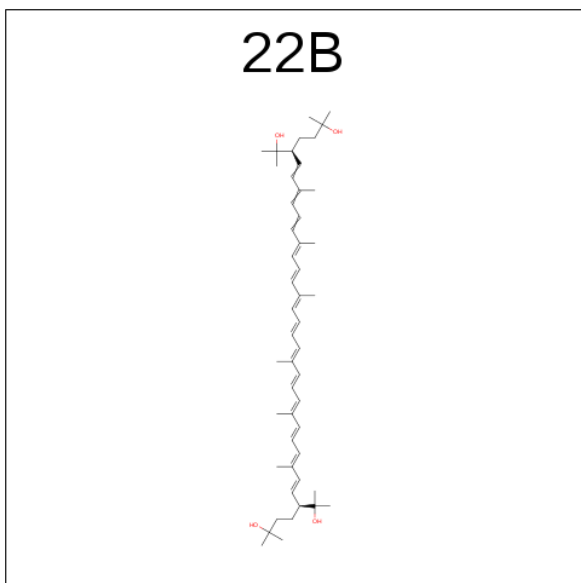
- Molecule 3 is RETINAL (three-letter code: RET) (formula: C<sub>20</sub>H<sub>28</sub>O).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C 20 20	0	0
3	B	1	Total C 20 20	0	0
3	D	1	Total C 20 20	0	0
3	E	1	Total C 20 20	0	0

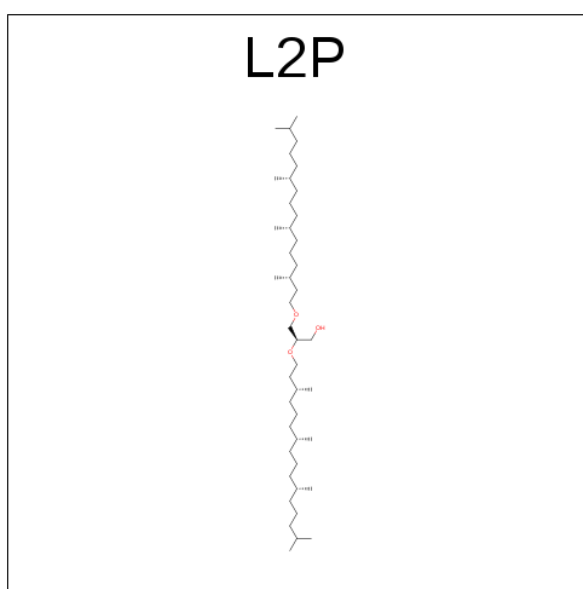
- Molecule 4 is BACTERIORUBERIN (three-letter code: 22B) (formula:  $C_{50}H_{76}O_4$ ).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			54	50	4		
4	B	1	Total	C	O	0	0
			54	50	4		
4	D	1	Total	C	O	0	0
			54	50	4		
4	E	1	Total	C	O	0	0
			54	50	4		

- Molecule 5 is 2,3-DI-PHYTANYL-GLYCEROL (three-letter code: L2P) (formula:  $C_{43}H_{88}O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			46	43	3		
5	B	1	Total	C	O	0	0
			46	43	3		
5	D	1	Total	C	O	0	0
			46	43	3		
5	E	1	Total	C	O	0	0
			46	43	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	25	Total	O	0	0
			25	25		

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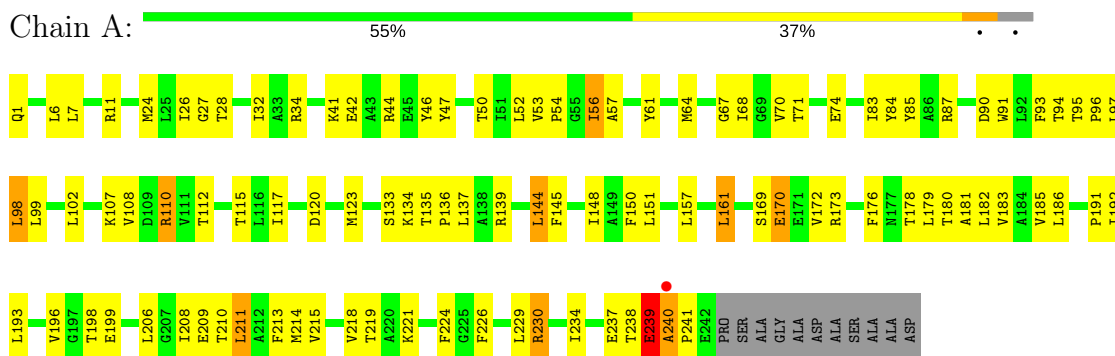
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	26	Total 26	O 26	0	0
6	D	22	Total 22	O 22	0	0
6	E	28	Total 28	O 28	0	0



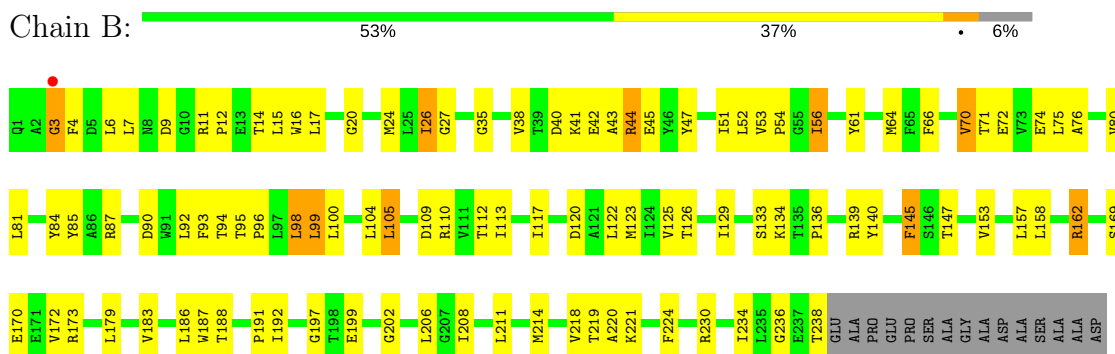
### 3 Residue-property plots

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.

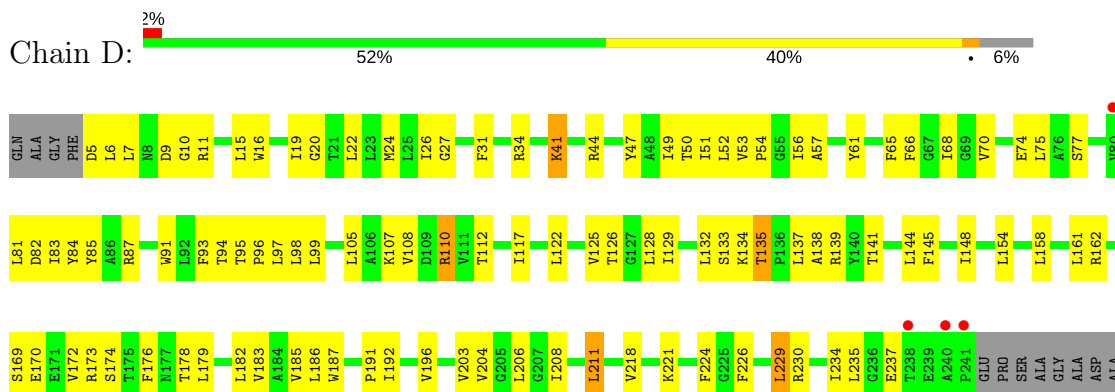
#### • Molecule 1: Archaerhodopsin-2



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SER

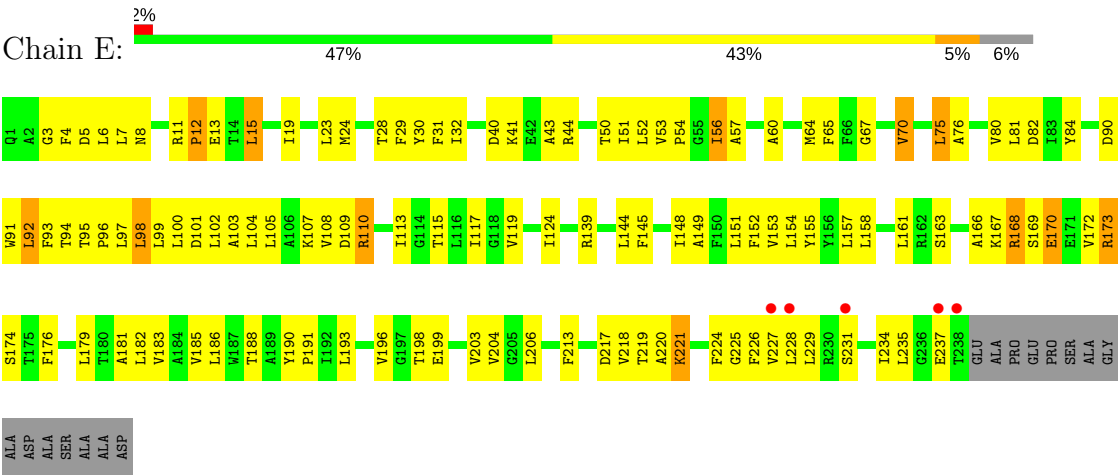
ALA

ALA

ALA

ASP

• Molecule 1: Archaerhodopsin-2





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	108.83Å 108.83Å 220.71Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	15.00 – 2.50 54.42 – 2.50	Depositor EDS
% Data completeness (in resolution range)	93.6 (15.00-2.50) 93.6 (54.42-2.50)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.71 (at 2.51Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.201 , 0.232 0.203 , 0.233	Depositor DCC
$R_{free}$ test set	4860 reflections (10.22%)	DCC
Wilson B-factor (Å <sup>2</sup> )	39.1	Xtriage
Anisotropy	0.255	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 73.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.44$ , $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.156 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	8087	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.75% of the height of the origin peak. No significant pseudotranslation is detected.*

<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: GLC, 22B, RET, L2P, GAL, MAN

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	0.38	0/1911	0.62	0/2612
1	B	0.36	0/1880	0.62	0/2569
1	D	0.34	0/1872	0.56	0/2560
1	E	0.31	0/1880	0.53	0/2569
All	All	0.35	0/7543	0.58	0/10310

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	1868	0	1929	99	0
1	B	1838	0	1904	79	0
1	D	1830	0	1894	87	0
1	E	1838	0	1905	91	0
2	A	33	0	28	0	0
2	B	33	0	28	0	0
2	D	66	0	56	3	0
3	A	20	0	27	2	0
3	B	20	0	27	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	20	0	27	0	0
3	E	20	0	27	2	0
4	A	54	0	76	3	0
4	B	54	0	76	2	0
4	D	54	0	76	2	0
4	E	54	0	76	0	0
5	A	46	0	87	5	0
5	B	46	0	87	1	0
5	D	46	0	87	3	0
5	E	46	0	87	2	0
6	A	25	0	0	3	0
6	B	26	0	0	1	0
6	D	22	0	0	2	0
6	E	28	0	0	3	0
All	All	8087	0	8504	353	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 21.

All (353) 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:135:THR:HG22	1:D:138:ALA:H	1.38	0.88
1:B:208:ILE:H	1:B:208:ILE:HD12	1.42	0.85
1:D:53:VAL:HG12	1:D:221:LYS:HG3	1.58	0.84
1:B:52:LEU:O	1:B:56:ILE:HG22	1.79	0.82
1:A:53:VAL:HG12	1:A:221:LYS:HG3	1.61	0.82
1:B:100:LEU:O	1:B:104:LEU:HD13	1.82	0.80
1:E:186:LEU:HD23	1:E:220:ALA:HB2	1.67	0.77
1:D:133:SER:HB2	1:D:139:ARG:HG2	1.67	0.77
1:A:170:GLU:HA	1:A:173:ARG:HG2	1.68	0.76
1:A:240:ALA:H	1:A:241:PRO:HD2	1.51	0.76
1:E:203:VAL:HG23	1:E:204:VAL:HG23	1.69	0.75
1:D:112:THR:HA	4:D:270:22B:H241	1.71	0.71
1:A:230:ARG:HG2	1:A:230:ARG:HH11	1.56	0.70
1:E:170:GLU:HA	1:E:173:ARG:HG2	1.73	0.70
1:A:108:VAL:HG11	1:A:161:LEU:HD22	1.74	0.70
1:E:53:VAL:HG21	1:E:98:LEU:HD13	1.73	0.69
1:E:166:ALA:HB1	1:E:173:ARG:HH11	1.55	0.69
1:A:226:PHE:O	1:A:230:ARG:HB2	1.92	0.69
1:A:173:ARG:HD2	6:A:296:HOH:O	1.94	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:ARG:O	1:A:90:ASP:HB3	1.93	0.68
1:E:41:LYS:HA	1:E:44:ARG:HD3	1.76	0.67
1:D:6:LEU:HD11	1:D:15:LEU:HA	1.76	0.67
1:E:52:LEU:O	1:E:56:ILE:HG22	1.95	0.67
1:D:144:LEU:O	1:D:148:ILE:HG12	1.94	0.66
1:A:133:SER:HB2	1:A:139:ARG:HG2	1.77	0.66
1:D:95:THR:OG1	1:D:96:PRO:HD3	1.95	0.66
1:D:84:TYR:HE2	1:D:206:LEU:HD21	1.60	0.66
1:B:15:LEU:HD13	1:B:15:LEU:O	1.96	0.65
1:A:112:THR:HA	4:A:270:22B:H241	1.77	0.65
1:B:188:THR:O	1:B:192:ILE:HD13	1.96	0.65
1:A:178:THR:HG22	1:A:178:THR:O	1.95	0.65
1:A:95:THR:OG1	1:A:96:PRO:HD3	1.95	0.65
1:E:91:TRP:HA	1:E:94:THR:OG1	1.96	0.65
1:D:122:LEU:HD21	4:D:270:22B:H203	1.77	0.65
1:B:53:VAL:HG12	1:B:221:LYS:HG3	1.77	0.64
1:D:226:PHE:O	1:D:230:ARG:HB2	1.97	0.64
1:D:41:LYS:HA	1:D:44:ARG:NH1	2.12	0.64
1:B:14:THR:HA	1:B:17:LEU:HD12	1.78	0.64
1:A:191:PRO:HB3	3:A:260:RET:H183	1.79	0.64
1:D:208:ILE:H	1:D:208:ILE:HD12	1.62	0.64
1:A:107:LYS:HE3	1:A:239:GLU:H	1.61	0.64
1:A:52:LEU:O	1:A:56:ILE:HG22	1.97	0.64
1:E:191:PRO:HB3	3:E:260:RET:H183	1.80	0.64
1:E:24:MET:HE1	1:E:221:LYS:HB3	1.79	0.64
1:A:144:LEU:O	1:A:148:ILE:HG12	1.97	0.63
1:E:168:ARG:HH21	1:E:237:GLU:HG2	1.63	0.63
1:E:40:ASP:HB3	1:E:43:ALA:HB3	1.80	0.63
1:D:34:ARG:HH22	1:D:230:ARG:HD2	1.64	0.63
1:B:120:ASP:O	1:B:123:MET:HB3	1.99	0.62
1:A:214:MET:O	1:A:218:VAL:HG12	1.99	0.62
1:A:110:ARG:NE	1:A:110:ARG:H	1.98	0.62
1:E:57:ALA:HA	1:E:90:ASP:OD2	1.99	0.62
1:A:240:ALA:N	1:A:241:PRO:HD2	2.15	0.61
1:B:208:ILE:HD12	1:B:208:ILE:N	2.15	0.61
1:E:99:LEU:HD12	1:E:102:LEU:HD12	1.80	0.61
1:B:95:THR:OG1	1:B:96:PRO:HD3	2.01	0.61
1:D:110:ARG:HD2	1:D:110:ARG:H	1.65	0.61
1:E:139:ARG:HH21	1:E:199:GLU:HA	1.64	0.61
1:B:26:ILE:HG13	1:B:27:GLY:N	2.15	0.60
1:A:182:LEU:O	1:A:186:LEU:HB2	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:LYS:HZ2	1:A:238:THR:HG23	1.65	0.60
1:B:40:ASP:HB3	1:B:43:ALA:HB3	1.82	0.60
2:D:286:GAL:H3	1:E:67:GLY:HA3	1.84	0.60
1:D:52:LEU:O	1:D:56:ILE:HG22	2.01	0.60
1:B:70:VAL:HG22	1:B:84:TYR:CE1	2.36	0.60
1:A:157:LEU:HD11	1:A:180:THR:HG23	1.83	0.60
1:A:206:LEU:HD23	1:A:206:LEU:O	2.01	0.60
1:E:96:PRO:HB3	1:E:117:ILE:HG13	1.82	0.60
1:A:41:LYS:HA	1:A:44:ARG:CD	2.32	0.59
1:E:107:LYS:HB2	1:E:168:ARG:HD2	1.84	0.59
1:E:169:SER:HB3	1:E:172:VAL:HG23	1.83	0.59
1:A:26:ILE:HG13	1:A:27:GLY:N	2.17	0.59
1:E:60:ALA:O	1:E:64:MET:HG3	2.03	0.59
1:E:102:LEU:HD13	1:E:157:LEU:HD11	1.85	0.58
1:A:115:THR:HG23	4:A:270:22B:H183	1.86	0.58
1:E:229:LEU:HA	1:E:234:ILE:HD13	1.85	0.58
1:E:95:THR:OG1	1:E:96:PRO:HD3	2.03	0.58
1:A:120:ASP:O	1:A:123:MET:HB3	2.04	0.58
1:B:136:PRO:O	1:B:140:TYR:HD1	1.86	0.58
1:A:170:GLU:HA	1:A:173:ARG:NE	2.19	0.57
1:A:198:THR:HG23	1:A:209:GLU:OE1	2.04	0.57
1:E:92:LEU:HD22	1:E:124:ILE:HB	1.87	0.57
1:D:66:PHE:O	2:D:283:GAL:H5	2.04	0.57
1:D:93:PHE:C	1:D:96:PRO:HD2	2.25	0.57
1:B:53:VAL:CG1	1:B:221:LYS:HG3	2.35	0.56
1:B:153:VAL:O	1:B:157:LEU:HB2	2.05	0.56
1:B:61:TYR:OH	1:B:214:MET:HA	2.06	0.56
1:D:70:VAL:HG22	1:D:84:TYR:CD1	2.41	0.56
1:A:135:THR:HG22	1:A:137:LEU:H	1.70	0.56
1:D:49:ILE:HG23	1:D:97:LEU:HD22	1.88	0.56
1:E:113:ILE:HG22	1:E:117:ILE:HD13	1.88	0.55
1:A:211:LEU:O	1:A:215:VAL:HG23	2.07	0.55
1:A:53:VAL:CG1	1:A:221:LYS:HG3	2.36	0.55
1:A:67:GLY:O	1:A:70:VAL:HG12	2.06	0.55
1:A:161:LEU:HB3	1:A:176:PHE:HZ	1.71	0.55
1:A:107:LYS:HG3	1:A:239:GLU:HB2	1.88	0.55
1:B:24:MET:CE	1:B:54:PRO:HA	2.37	0.55
1:D:96:PRO:HB3	1:D:117:ILE:HG13	1.89	0.55
1:B:187:TRP:O	1:B:191:PRO:HD2	2.07	0.55
1:A:34:ARG:HH12	1:A:230:ARG:HE	1.55	0.54
1:D:183:VAL:HG13	1:D:187:TRP:CE2	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:LEU:HD23	1:B:76:ALA:N	2.23	0.54
1:E:3:GLY:HA3	1:E:12:PRO:O	2.07	0.54
1:E:168:ARG:NH2	1:E:237:GLU:HG2	2.23	0.54
1:D:192:ILE:O	1:D:196:VAL:HG22	2.08	0.54
1:B:92:LEU:HD23	1:B:93:PHE:CE2	2.43	0.54
1:A:133:SER:HB2	1:A:139:ARG:CG	2.38	0.54
1:D:185:VAL:HG13	1:D:186:LEU:HD12	1.90	0.54
1:D:34:ARG:HH22	1:D:230:ARG:CD	2.20	0.54
1:E:110:ARG:H	1:E:110:ARG:HD2	1.72	0.54
1:D:126:THR:HG21	1:D:145:PHE:CD2	2.42	0.53
1:D:126:THR:HG21	1:D:145:PHE:CE2	2.44	0.53
1:D:7:LEU:HD11	1:D:65:PHE:CZ	2.43	0.53
1:E:82:ASP:HA	6:E:307:HOH:O	2.09	0.53
1:A:161:LEU:HB3	1:A:176:PHE:CZ	2.42	0.53
1:D:182:LEU:O	1:D:186:LEU:HD13	2.09	0.53
1:E:6:LEU:HD11	1:E:15:LEU:HD22	1.90	0.53
1:E:40:ASP:O	1:E:44:ARG:HG2	2.09	0.53
1:A:91:TRP:HA	1:A:94:THR:OG1	2.09	0.53
1:A:50:THR:HG21	1:A:229:LEU:HD21	1.90	0.53
1:D:74:GLU:O	1:D:134:LYS:HG3	2.09	0.53
1:B:113:ILE:O	1:B:117:ILE:HD13	2.09	0.52
1:D:75:LEU:HA	1:D:134:LYS:HG3	1.90	0.52
1:E:190:TYR:N	1:E:191:PRO:HD2	2.25	0.52
1:A:170:GLU:HA	1:A:173:ARG:CG	2.38	0.52
1:A:170:GLU:HA	1:A:173:ARG:HE	1.74	0.52
1:D:128:LEU:HD11	1:D:132:LEU:HD11	1.92	0.52
1:D:87:ARG:HG2	1:D:87:ARG:HH11	1.74	0.52
1:B:208:ILE:H	1:B:208:ILE:CD1	2.14	0.52
1:D:47:TYR:O	1:D:51:ILE:HG13	2.10	0.52
1:E:11:ARG:O	1:E:13:GLU:N	2.43	0.52
1:A:192:ILE:O	1:A:196:VAL:HG22	2.10	0.52
1:B:94:THR:O	1:B:98:LEU:HD22	2.10	0.52
1:A:24:MET:HE1	1:A:218:VAL:HA	1.90	0.51
1:B:133:SER:HB2	1:B:139:ARG:HG2	1.93	0.51
1:D:34:ARG:HH22	1:D:230:ARG:HE	1.58	0.51
1:B:74:GLU:HG2	1:B:80:VAL:HG22	1.93	0.51
1:D:125:VAL:O	1:D:129:ILE:HG13	2.10	0.51
1:D:26:ILE:HG13	1:D:27:GLY:N	2.25	0.51
1:A:70:VAL:HB	1:A:84:TYR:CE1	2.45	0.51
1:B:53:VAL:HB	1:B:54:PRO:CD	2.40	0.51
1:E:19:ILE:O	1:E:23:LEU:HG	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:GLY:O	1:B:12:PRO:HB3	2.11	0.51
1:D:50:THR:O	1:D:54:PRO:HD2	2.11	0.51
1:D:68:ILE:C	1:D:70:VAL:H	2.14	0.51
1:D:170:GLU:HG2	1:D:173:ARG:NH1	2.26	0.51
1:D:34:ARG:HH22	1:D:230:ARG:NE	2.09	0.51
1:A:1:GLN:HG3	1:E:30:TYR:OH	2.11	0.51
1:A:179:LEU:HB3	1:A:224:PHE:CE1	2.46	0.51
1:A:230:ARG:HG2	1:A:230:ARG:NH1	2.25	0.51
1:B:109:ASP:O	1:B:113:ILE:HG13	2.11	0.51
1:A:123:MET:SD	1:A:150:PHE:HB2	2.51	0.50
1:E:97:LEU:O	1:E:100:LEU:HB3	2.11	0.50
1:B:169:SER:OG	1:B:172:VAL:HG23	2.11	0.50
1:B:41:LYS:HG3	1:B:44:ARG:NH1	2.26	0.50
1:B:126:THR:HG21	1:B:145:PHE:CE2	2.45	0.50
1:A:102:LEU:HD11	1:A:183:VAL:HG21	1.94	0.50
1:A:24:MET:CE	1:A:218:VAL:HA	2.42	0.50
1:B:122:LEU:HD21	4:B:270:22B:H203	1.94	0.50
1:B:93:PHE:C	1:B:96:PRO:HD2	2.31	0.50
1:D:169:SER:OG	1:D:172:VAL:HG23	2.11	0.50
1:D:22:LEU:O	1:D:26:ILE:HG23	2.11	0.50
1:E:11:ARG:C	1:E:13:GLU:H	2.14	0.50
1:B:80:VAL:C	1:B:81:LEU:HD12	2.32	0.50
1:A:185:VAL:HG21	1:E:19:ILE:HD13	1.94	0.50
1:A:47:TYR:HA	1:A:229:LEU:HD11	1.93	0.50
1:B:24:MET:HE3	1:B:54:PRO:HA	1.94	0.50
1:D:183:VAL:HG13	1:D:187:TRP:CD1	2.47	0.50
1:D:187:TRP:O	1:D:191:PRO:HD2	2.12	0.49
1:E:217:ASP:O	1:E:221:LYS:HB2	2.12	0.49
1:A:181:ALA:O	1:A:185:VAL:HG12	2.11	0.49
1:A:191:PRO:HB3	3:A:260:RET:C18	2.42	0.49
1:B:24:MET:CE	1:B:221:LYS:HG2	2.42	0.49
1:D:179:LEU:HB3	1:D:224:PHE:CE1	2.47	0.49
1:D:230:ARG:HG2	1:D:230:ARG:HH11	1.77	0.49
1:A:218:VAL:HG13	1:A:219:THR:N	2.28	0.49
1:A:56:ILE:HD12	5:A:280:L2P:H53	1.93	0.49
1:B:197:GLY:O	1:B:202:GLY:HA2	2.13	0.49
1:E:53:VAL:HG12	1:E:221:LYS:HD2	1.95	0.49
1:E:31:PHE:CZ	1:E:225:GLY:HA3	2.47	0.49
1:A:107:LYS:NZ	1:A:238:THR:HA	2.28	0.49
1:D:183:VAL:HG13	1:D:187:TRP:NE1	2.28	0.49
1:E:154:LEU:O	1:E:158:LEU:HG	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:83:ILE:HA	6:D:293:HOH:O	2.13	0.48
1:E:144:LEU:O	1:E:148:ILE:HG12	2.13	0.48
1:E:50:THR:O	1:E:54:PRO:HD2	2.13	0.48
1:E:105:LEU:HD23	1:E:176:PHE:HD1	1.78	0.48
1:E:23:LEU:HD13	1:E:218:VAL:HG21	1.95	0.48
1:B:218:VAL:HG13	1:B:219:THR:N	2.29	0.48
1:D:53:VAL:CG1	1:D:221:LYS:HG3	2.38	0.48
1:E:7:LEU:N	1:E:7:LEU:HD12	2.29	0.48
1:D:24:MET:HE1	1:D:218:VAL:HA	1.95	0.48
1:E:115:THR:O	1:E:119:VAL:HG23	2.14	0.48
1:B:87:ARG:O	1:B:90:ASP:HB3	2.14	0.48
1:E:149:ALA:O	1:E:153:VAL:HG23	2.13	0.48
1:E:182:LEU:O	1:E:186:LEU:HB2	2.13	0.48
1:A:28:THR:O	1:A:32:ILE:HG13	2.14	0.47
1:E:67:GLY:O	1:E:70:VAL:HG23	2.14	0.47
1:D:108:VAL:HG11	1:D:161:LEU:HD13	1.95	0.47
1:D:185:VAL:HG13	1:D:186:LEU:N	2.29	0.47
1:D:68:ILE:HG12	2:D:283:GAL:O2	2.14	0.47
1:E:173:ARG:HG3	1:E:174:SER:N	2.28	0.47
1:B:199:GLU:OE1	1:B:199:GLU:N	2.47	0.47
1:B:11:ARG:NH2	1:B:64:MET:O	2.47	0.47
1:D:5:ASP:CG	1:D:10:GLY:H	2.17	0.47
1:A:11:ARG:NH2	1:A:64:MET:O	2.47	0.47
1:B:125:VAL:O	1:B:129:ILE:HG13	2.15	0.47
1:A:1:GLN:HB3	1:A:208:ILE:HD11	1.97	0.47
1:B:183:VAL:HG13	1:B:187:TRP:CE2	2.50	0.47
1:E:24:MET:CE	1:E:218:VAL:HA	2.45	0.47
1:A:71:THR:HG22	1:A:85:TYR:HD2	1.79	0.47
1:B:179:LEU:HD22	1:B:224:PHE:CD1	2.50	0.47
1:E:28:THR:O	1:E:32:ILE:HG13	2.15	0.47
1:E:31:PHE:HB2	1:E:51:ILE:HG12	1.96	0.47
1:D:75:LEU:HD23	1:D:77:SER:OG	2.14	0.47
1:B:126:THR:HG21	1:B:145:PHE:CD2	2.50	0.47
1:E:24:MET:HE3	1:E:218:VAL:HA	1.97	0.47
1:B:186:LEU:HD23	1:B:220:ALA:HB2	1.96	0.46
1:D:158:LEU:O	1:D:162:ARG:HD3	2.16	0.46
1:A:50:THR:HG21	1:A:229:LEU:CD2	2.45	0.46
1:B:9:ASP:CG	1:B:11:ARG:HD3	2.35	0.46
1:A:34:ARG:HH12	1:A:230:ARG:NE	2.14	0.46
1:B:112:THR:HA	4:B:270:22B:H241	1.97	0.46
1:A:148:ILE:O	1:A:151:LEU:HB3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:203:VAL:HG23	1:D:204:VAL:HG23	1.97	0.46
1:E:188:THR:O	1:E:191:PRO:HG2	2.14	0.46
1:A:50:THR:O	1:A:54:PRO:HD2	2.16	0.46
1:A:96:PRO:HB3	1:A:117:ILE:HG13	1.98	0.46
1:B:52:LEU:HD13	5:B:280:L2P:H593	1.96	0.46
1:D:15:LEU:O	1:D:19:ILE:HG13	2.15	0.46
1:E:103:ALA:HB1	1:E:113:ILE:HG12	1.98	0.46
1:B:179:LEU:HB3	1:B:224:PHE:HE1	1.81	0.46
1:D:70:VAL:HG22	1:D:84:TYR:CE1	2.51	0.46
1:B:24:MET:HE2	1:B:221:LYS:HG2	1.96	0.46
1:B:41:LYS:HG3	1:B:44:ARG:HH12	1.80	0.46
1:A:139:ARG:HH21	1:A:199:GLU:HA	1.81	0.46
1:A:169:SER:OG	1:A:172:VAL:HG23	2.16	0.46
1:A:179:LEU:O	1:A:183:VAL:HG23	2.16	0.46
1:B:211:LEU:HD11	1:D:185:VAL:CG2	2.46	0.46
1:D:16:TRP:CE3	1:D:211:LEU:HD12	2.51	0.46
1:E:186:LEU:HD21	1:E:219:THR:CG2	2.46	0.46
1:E:7:LEU:HD11	1:E:65:PHE:CZ	2.51	0.45
1:A:34:ARG:HH22	1:A:230:ARG:NE	2.14	0.45
1:B:6:LEU:HD11	1:B:15:LEU:HD23	1.99	0.45
1:E:50:THR:O	1:E:54:PRO:CD	2.65	0.45
1:E:75:LEU:HG	1:E:76:ALA:H	1.80	0.45
1:A:97:LEU:HD11	5:A:280:L2P:H592	1.98	0.45
1:B:35:GLY:HA2	1:B:38:VAL:HG23	1.98	0.45
1:D:91:TRP:HA	1:D:94:THR:OG1	2.17	0.45
1:B:9:ASP:OD1	1:B:11:ARG:HD3	2.18	0.44
1:B:183:VAL:HG13	1:B:187:TRP:CD2	2.51	0.44
1:D:56:ILE:HG13	5:D:280:L2P:H53	1.97	0.44
1:B:16:TRP:O	1:B:214:MET:HG3	2.17	0.44
1:B:61:TYR:HA	1:B:64:MET:HE3	1.99	0.44
1:D:9:ASP:HB3	1:D:11:ARG:HD3	1.99	0.44
1:D:137:LEU:O	1:D:141:THR:HG22	2.17	0.44
1:E:70:VAL:HG22	1:E:84:TYR:CE1	2.52	0.44
1:E:15:LEU:CD1	1:E:19:ILE:HD11	2.48	0.44
1:E:93:PHE:C	1:E:96:PRO:HD2	2.37	0.44
1:B:15:LEU:HD13	1:B:15:LEU:C	2.38	0.44
1:E:193:LEU:HD23	1:E:213:PHE:CE1	2.52	0.44
1:A:34:ARG:HH22	1:A:230:ARG:CD	2.31	0.44
1:A:74:GLU:O	1:A:134:LYS:HG3	2.17	0.44
1:D:135:THR:O	1:D:139:ARG:HG3	2.18	0.44
1:B:234:ILE:C	1:B:236:GLY:H	2.22	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:152:PHE:O	1:E:155:TYR:HB3	2.18	0.44
1:E:139:ARG:NH2	1:E:199:GLU:HA	2.32	0.44
1:E:56:ILE:HD12	5:E:280:L2P:H53	2.00	0.44
1:E:5:ASP:OD2	1:E:8:ASN:HA	2.16	0.44
1:B:38:VAL:HG21	1:B:47:TYR:CE1	2.53	0.44
1:A:115:THR:CG2	4:A:270:22B:H183	2.48	0.43
1:D:133:SER:HB2	1:D:139:ARG:CG	2.41	0.43
1:D:7:LEU:HD12	1:D:7:LEU:N	2.34	0.43
1:B:238:THR:HG22	6:B:304:HOH:O	2.17	0.43
1:A:61:TYR:OH	1:A:214:MET:HA	2.19	0.43
1:A:57:ALA:HA	1:A:90:ASP:OD2	2.18	0.43
1:B:71:THR:HG23	1:B:85:TYR:CD2	2.53	0.43
1:A:206:LEU:HD23	1:A:210:THR:OG1	2.17	0.43
1:E:166:ALA:HB1	1:E:173:ARG:NH1	2.29	0.43
1:E:181:ALA:O	1:E:185:VAL:HG12	2.19	0.43
1:E:179:LEU:O	1:E:183:VAL:HG23	2.18	0.43
1:E:198:THR:HA	6:E:301:HOH:O	2.18	0.43
1:D:173:ARG:HG3	1:D:174:SER:N	2.33	0.43
1:A:56:ILE:HG23	1:A:94:THR:CG2	2.49	0.43
1:D:75:LEU:CA	1:D:134:LYS:HG3	2.48	0.43
1:A:234:ILE:HD11	6:A:303:HOH:O	2.19	0.43
1:A:42:GLU:HG3	1:A:46:TYR:HE2	1.84	0.43
1:D:170:GLU:HG3	6:E:293:HOH:O	2.19	0.43
1:B:211:LEU:HD23	1:B:211:LEU:C	2.40	0.42
1:D:87:ARG:HG2	1:D:87:ARG:NH1	2.34	0.42
1:B:147:THR:HA	3:B:260:RET:H182	2.02	0.42
1:B:47:TYR:O	1:B:51:ILE:HG13	2.19	0.42
1:E:100:LEU:O	1:E:104:LEU:HD13	2.19	0.42
1:B:42:GLU:O	1:B:45:GLU:HB3	2.19	0.42
1:E:29:PHE:HA	1:E:32:ILE:HD12	2.01	0.42
1:A:6:LEU:O	1:A:7:LEU:HD23	2.19	0.42
1:A:93:PHE:C	1:A:96:PRO:HD2	2.40	0.42
1:D:105:LEU:HD23	1:D:176:PHE:HA	2.00	0.42
1:D:31:PHE:CB	1:D:51:ILE:HG12	2.49	0.42
1:A:135:THR:HG23	1:A:136:PRO:HD2	2.01	0.42
1:A:85:TYR:CE1	5:A:280:L2P:H43	2.54	0.42
1:A:85:TYR:HE1	5:A:280:L2P:H43	1.85	0.42
1:E:166:ALA:CB	1:E:173:ARG:HE	2.32	0.42
1:A:50:THR:CG2	1:A:229:LEU:HD21	2.49	0.42
1:B:74:GLU:O	1:B:134:LYS:HE3	2.19	0.42
1:D:144:LEU:HD23	1:D:144:LEU:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:85:TYR:OH	5:D:280:L2P:H43	2.20	0.42
1:E:196:VAL:HB	1:E:203:VAL:HG22	2.01	0.42
1:A:107:LYS:HZ1	1:A:238:THR:HA	1.85	0.41
1:E:179:LEU:O	1:E:182:LEU:HB3	2.20	0.41
1:B:95:THR:O	1:B:99:LEU:HD22	2.20	0.41
1:D:56:ILE:HG23	1:D:57:ALA:N	2.36	0.41
1:A:135:THR:HG22	1:A:137:LEU:N	2.36	0.41
1:E:148:ILE:O	1:E:151:LEU:HB3	2.19	0.41
1:A:170:GLU:CA	1:A:173:ARG:HG2	2.43	0.41
1:A:83:ILE:HG23	6:A:291:HOH:O	2.21	0.41
1:B:20:GLY:HA3	1:B:61:TYR:CE2	2.56	0.41
1:B:4:PHE:CE2	1:D:178:THR:HG23	2.56	0.41
1:E:101:ASP:OD1	1:E:228:LEU:HD21	2.20	0.41
1:D:47:TYR:CE2	1:D:229:LEU:HG	2.56	0.41
1:E:163:SER:O	1:E:167:LYS:HG3	2.21	0.41
1:E:179:LEU:HD21	1:E:227:VAL:HB	2.03	0.41
1:D:20:GLY:HA3	1:D:61:TYR:CD2	2.55	0.41
1:D:31:PHE:HB3	1:D:51:ILE:HG12	2.03	0.41
1:E:234:ILE:HG13	1:E:235:LEU:N	2.36	0.41
1:E:80:VAL:C	1:E:81:LEU:HD12	2.41	0.41
1:A:185:VAL:HG21	1:E:19:ILE:CD1	2.50	0.41
1:A:93:PHE:O	1:A:97:LEU:HG	2.21	0.41
1:A:52:LEU:HD13	5:A:280:L2P:H593	2.02	0.41
5:D:280:L2P:H12	5:D:280:L2P:H411	1.89	0.41
1:E:221:LYS:O	1:E:224:PHE:HB3	2.21	0.41
1:A:193:LEU:HD23	1:A:213:PHE:CE1	2.55	0.41
1:B:211:LEU:HD11	1:D:185:VAL:HG21	2.02	0.41
1:E:108:VAL:HG11	1:E:161:LEU:HD13	2.02	0.41
1:A:144:LEU:HD22	1:A:148:ILE:HD11	2.03	0.41
1:A:68:ILE:H	1:A:68:ILE:HG13	1.54	0.41
1:D:24:MET:HE2	1:D:24:MET:HA	2.03	0.41
1:E:98:LEU:O	1:E:101:ASP:HB2	2.21	0.41
1:B:71:THR:HG23	1:B:85:TYR:CE2	2.55	0.40
1:D:53:VAL:HB	1:D:54:PRO:CD	2.51	0.40
1:E:196:VAL:O	1:E:203:VAL:HG22	2.20	0.40
1:E:98:LEU:HD12	1:E:98:LEU:HA	1.87	0.40
1:A:34:ARG:NH1	1:A:230:ARG:HE	2.16	0.40
1:B:170:GLU:HA	1:B:173:ARG:HG2	2.04	0.40
1:D:138:ALA:HA	1:D:141:THR:HG22	2.03	0.40
1:D:234:ILE:HD11	6:D:307:HOH:O	2.21	0.40
1:B:158:LEU:HA	1:B:162:ARG:HG3	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:7:LEU:HD21	1:B:66:PHE:HA	2.03	0.40
1:A:98:LEU:HD13	1:A:98:LEU:HA	1.93	0.40
1:D:154:LEU:O	1:D:158:LEU:HG	2.21	0.40
3:E:260:RET:H7	3:E:260:RET:H181	1.89	0.40
5:E:280:L2P:H12	5:E:280:L2P:H411	1.91	0.40
1:B:105:LEU:HD12	1:B:105:LEU:HA	1.96	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	240/253 (95%)	226 (94%)	11 (5%)	3 (1%)	14	25
1	B	236/253 (93%)	215 (91%)	20 (8%)	1 (0%)	38	59
1	D	235/253 (93%)	219 (93%)	14 (6%)	2 (1%)	20	36
1	E	236/253 (93%)	217 (92%)	14 (6%)	5 (2%)	8	13
All	All	947/1012 (94%)	877 (93%)	59 (6%)	11 (1%)	15	27

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	240	ALA
1	D	237	GLU
1	E	231	SER
1	D	82	ASP
1	E	109	ASP
1	A	237	GLU
1	A	239	GLU
1	E	168	ARG
1	E	75	LEU

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Mol	Chain	Res	Type
1	E	12	PRO
1	B	3	GLY

### 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	191/196 (97%)	180 (94%)	11 (6%)	23	43
1	B	188/196 (96%)	175 (93%)	13 (7%)	18	34
1	D	188/196 (96%)	178 (95%)	10 (5%)	26	48
1	E	188/196 (96%)	175 (93%)	13 (7%)	18	34
All	All	755/784 (96%)	708 (94%)	47 (6%)	21	39

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	ILE
1	A	98	LEU
1	A	99	LEU
1	A	110	ARG
1	A	144	LEU
1	A	145	PHE
1	A	161	LEU
1	A	170	GLU
1	A	211	LEU
1	A	230	ARG
1	A	239	GLU
1	B	26	ILE
1	B	44	ARG
1	B	56	ILE
1	B	70	VAL
1	B	72	GLU
1	B	98	LEU
1	B	99	LEU
1	B	105	LEU

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Mol	Chain	Res	Type
1	B	110	ARG
1	B	145	PHE
1	B	162	ARG
1	B	206	LEU
1	B	230	ARG
1	D	41	LYS
1	D	81	LEU
1	D	98	LEU
1	D	99	LEU
1	D	107	LYS
1	D	110	ARG
1	D	135	THR
1	D	211	LEU
1	D	229	LEU
1	D	235	LEU
1	E	4	PHE
1	E	15	LEU
1	E	56	ILE
1	E	70	VAL
1	E	92	LEU
1	E	98	LEU
1	E	110	ARG
1	E	145	PHE
1	E	170	GLU
1	E	173	ARG
1	E	206	LEU
1	E	221	LYS
1	E	226	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	177	ASN
1	B	177	ASN
1	D	177	ASN
1	E	8	ASN
1	E	177	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 ⓘ

12 carbohydrates 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	GLC	A	281	2,5	11,11,12	2.02	3 (27%)	13,15,17	2.05	6 (46%)
2	MAN	A	282	2	11,11,12	2.49	7 (63%)	13,15,17	0.88	0
2	GAL	A	283	2	11,11,12	1.88	3 (27%)	13,15,17	1.05	0
2	GLC	B	281	2,5	11,11,12	1.26	1 (9%)	13,15,17	2.15	5 (38%)
2	MAN	B	282	2	11,11,12	2.31	5 (45%)	13,15,17	1.48	3 (23%)
2	GAL	B	283	2	11,11,12	1.97	3 (27%)	13,15,17	2.34	2 (15%)
2	GLC	D	281	2,5	11,11,12	2.01	2 (18%)	13,15,17	1.45	1 (7%)
2	MAN	D	282	2	11,11,12	2.66	7 (63%)	13,15,17	0.84	0
2	GAL	D	283	2	11,11,12	1.86	4 (36%)	13,15,17	1.05	1 (7%)
2	GLC	D	284	2,5	11,11,12	1.71	3 (27%)	13,15,17	1.40	1 (7%)
2	MAN	D	285	2	11,11,12	2.48	7 (63%)	13,15,17	0.79	0
2	GAL	D	286	2	11,11,12	1.56	3 (27%)	13,15,17	1.01	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	GLC	A	281	2,5	-	0/2/19/22	0/1/1/1
2	MAN	A	282	2	-	0/2/19/22	0/1/1/1
2	GAL	A	283	2	-	0/2/19/22	0/1/1/1
2	GLC	B	281	2,5	-	0/2/19/22	0/1/1/1
2	MAN	B	282	2	-	0/2/19/22	0/1/1/1
2	GAL	B	283	2	-	0/2/19/22	0/1/1/1
2	GLC	D	281	2,5	-	0/2/19/22	0/1/1/1
2	MAN	D	282	2	-	0/2/19/22	0/1/1/1
2	GAL	D	283	2	-	0/2/19/22	0/1/1/1
2	GLC	D	284	2,5	-	0/2/19/22	0/1/1/1
2	MAN	D	285	2	-	0/2/19/22	0/1/1/1
2	GAL	D	286	2	-	0/2/19/22	0/1/1/1

All (48) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	283	GAL	O5-C5	2.05	1.47	1.43
2	D	286	GAL	O5-C1	2.17	1.47	1.43
2	D	283	GAL	C4-C5	2.24	1.57	1.53
2	D	286	GAL	C4-C5	2.24	1.57	1.53
2	B	282	MAN	C4-C3	2.25	1.58	1.52
2	B	281	GLC	C1-C2	2.26	1.57	1.52
2	A	282	MAN	C4-C5	2.28	1.57	1.53
2	A	283	GAL	C4-C5	2.29	1.57	1.53
2	B	283	GAL	C4-C3	2.29	1.58	1.52
2	D	285	MAN	C6-C5	2.40	1.60	1.51
2	D	282	MAN	C4-C5	2.42	1.58	1.53
2	A	281	GLC	C1-C2	2.49	1.58	1.52
2	D	284	GLC	C1-C2	2.50	1.58	1.52
2	D	285	MAN	C4-C5	2.52	1.58	1.53
2	B	282	MAN	O5-C1	2.54	1.47	1.43
2	D	284	GLC	C4-C5	2.61	1.58	1.53
2	A	282	MAN	C6-C5	2.61	1.60	1.51
2	A	282	MAN	O5-C5	2.66	1.49	1.43
2	A	281	GLC	C4-C5	2.68	1.58	1.53
2	A	282	MAN	C4-C3	2.72	1.59	1.52
2	B	283	GAL	C4-C5	2.72	1.58	1.53
2	D	286	GAL	C2-C3	2.75	1.56	1.52
2	D	282	MAN	C6-C5	2.83	1.61	1.51
2	A	282	MAN	C2-C3	2.87	1.56	1.52
2	D	282	MAN	C1-C2	2.88	1.59	1.52
2	A	283	GAL	O5-C1	2.89	1.48	1.43
2	D	282	MAN	C4-C3	2.92	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	283	GAL	O5-C1	3.01	1.48	1.43
2	D	285	MAN	C4-C3	3.01	1.60	1.52
2	D	285	MAN	O5-C5	3.02	1.49	1.43
2	A	282	MAN	C1-C2	3.08	1.59	1.52
2	D	285	MAN	C1-C2	3.08	1.59	1.52
2	D	282	MAN	O5-C5	3.12	1.50	1.43
2	D	281	GLC	C1-C2	3.21	1.59	1.52
2	B	282	MAN	C4-C5	3.26	1.60	1.53
2	D	282	MAN	C2-C3	3.30	1.57	1.52
2	D	283	GAL	C2-C3	3.31	1.57	1.52
2	D	284	GLC	C2-C3	3.33	1.57	1.52
2	D	285	MAN	C2-C3	3.34	1.57	1.52
2	B	282	MAN	C1-C2	3.59	1.60	1.52
2	A	283	GAL	C2-C3	3.64	1.57	1.52
2	D	285	MAN	O5-C1	3.79	1.49	1.43
2	B	282	MAN	C2-C3	4.08	1.58	1.52
2	D	281	GLC	C2-C3	4.09	1.58	1.52
2	A	282	MAN	O5-C1	4.57	1.51	1.43
2	A	281	GLC	C2-C3	4.60	1.58	1.52
2	D	282	MAN	O5-C1	4.86	1.51	1.43
2	B	283	GAL	C2-C3	5.07	1.59	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	283	GAL	C1-C2-C3	-6.95	100.84	109.65
2	B	281	GLC	O2-C2-C3	-3.93	102.46	110.17
2	B	282	MAN	O6-C6-C5	-3.41	99.88	111.34
2	B	281	GLC	C6-C5-C4	-3.23	105.44	113.00
2	B	283	GAL	C1-O5-C5	-2.87	108.21	112.17
2	B	281	GLC	C3-C4-C5	-2.74	105.39	110.22
2	B	281	GLC	O4-C4-C3	-2.59	104.73	110.36
2	B	282	MAN	C2-C3-C4	-2.45	106.59	110.88
2	A	281	GLC	C3-C4-C5	-2.45	105.90	110.22
2	D	284	GLC	C3-C4-C5	-2.34	106.09	110.22
2	D	283	GAL	C1-C2-C3	-2.16	106.91	109.65
2	B	282	MAN	C6-C5-C4	-2.16	107.95	113.00
2	A	281	GLC	O2-C2-C1	-2.01	105.08	109.18
2	A	281	GLC	C6-C5-C4	2.13	117.99	113.00
2	A	281	GLC	C1-O5-C5	2.31	115.35	112.17
2	A	281	GLC	O6-C6-C5	2.95	121.27	111.34
2	B	281	GLC	O6-C6-C5	3.22	122.17	111.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	281	GLC	C1-O5-C5	3.87	117.50	112.17
2	A	281	GLC	O2-C2-C3	4.63	119.28	110.17

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	D	283	GAL	2	0
2	D	286	GAL	1	0

## 5.6 Ligand geometry [i](#)

12 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	RET	A	260	1	19,20,21	2.78	5 (26%)	27,27,28	2.12	15 (55%)
4	22B	A	270	-	51,53,53	2.23	10 (19%)	60,72,72	1.42	13 (21%)
5	L2P	A	280	2	45,45,45	1.26	5 (11%)	49,53,53	0.80	0
3	RET	B	260	1	19,20,21	2.91	6 (31%)	27,27,28	2.00	14 (51%)
4	22B	B	270	-	51,53,53	2.16	10 (19%)	60,72,72	1.53	14 (23%)
5	L2P	B	280	2	45,45,45	1.22	4 (8%)	49,53,53	0.83	0
3	RET	D	260	1	19,20,21	2.87	6 (31%)	27,27,28	2.07	14 (51%)
4	22B	D	270	-	51,53,53	2.21	10 (19%)	60,72,72	1.58	14 (23%)
5	L2P	D	280	2	45,45,45	1.27	4 (8%)	49,53,53	0.80	0
3	RET	E	260	1	19,20,21	3.01	7 (36%)	27,27,28	2.07	14 (51%)
4	22B	E	270	-	51,53,53	2.36	9 (17%)	60,72,72	1.54	14 (23%)
5	L2P	E	280	2	45,45,45	1.21	4 (8%)	49,53,53	0.86	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	RET	A	260	1	-	0/13/30/31	0/1/1/1
4	22B	A	270	-	-	0/65/65/65	0/0/0/0
5	L2P	A	280	2	-	0/51/51/51	0/0/0/0
3	RET	B	260	1	-	0/13/30/31	0/1/1/1
4	22B	B	270	-	-	0/65/65/65	0/0/0/0
5	L2P	B	280	2	-	0/51/51/51	0/0/0/0
3	RET	D	260	1	-	0/13/30/31	0/1/1/1
4	22B	D	270	-	-	0/65/65/65	0/0/0/0
5	L2P	D	280	2	-	0/51/51/51	0/0/0/0
3	RET	E	260	1	-	0/13/30/31	0/1/1/1
4	22B	E	270	-	-	0/65/65/65	0/0/0/0
5	L2P	E	280	2	-	0/51/51/51	0/0/0/0

All (80) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	E	280	L2P	C42-C41	2.00	1.57	1.50
5	A	280	L2P	C44-C43	2.00	1.59	1.52
5	A	280	L2P	O2-C41	2.02	1.48	1.43
3	D	260	RET	C10-C9	2.04	1.38	1.35
5	E	280	L2P	C3-C2	2.05	1.56	1.50
5	D	280	L2P	O2-C41	2.09	1.48	1.43
4	B	270	22B	C22-C21	2.11	1.58	1.53
4	B	270	22B	C21-C2	2.12	1.56	1.53
4	D	270	22B	C21-C2	2.12	1.56	1.53
5	B	280	L2P	O2-C41	2.12	1.48	1.43
4	D	270	22B	C22-C21	2.13	1.58	1.53
5	E	280	L2P	C19-C18	2.14	1.59	1.52
3	B	260	RET	C8-C7	2.19	1.39	1.33
3	E	260	RET	C8-C7	2.19	1.39	1.33
5	D	280	L2P	C42-C41	2.19	1.58	1.50
3	E	260	RET	C4-C5	2.20	1.55	1.51
5	B	280	L2P	C1-C2	2.26	1.57	1.50
4	A	270	22B	C21-C2	2.26	1.56	1.53
5	B	280	L2P	C3-C2	2.27	1.57	1.50
5	B	280	L2P	C42-C41	2.28	1.58	1.50
5	A	280	L2P	C42-C41	2.28	1.58	1.50
5	A	280	L2P	C1-C2	2.28	1.57	1.50
5	A	280	L2P	C3-C2	2.29	1.57	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	270	22B	C22-C21	2.35	1.58	1.53
5	E	280	L2P	C1-C2	2.35	1.57	1.50
5	D	280	L2P	C1-C2	2.38	1.57	1.50
5	D	280	L2P	C3-C2	2.50	1.57	1.50
3	B	260	RET	C18-C5	2.59	1.55	1.51
3	A	260	RET	C2-C1	2.60	1.60	1.54
3	B	260	RET	C2-C1	2.61	1.60	1.54
3	D	260	RET	C18-C5	2.70	1.55	1.51
3	E	260	RET	C2-C1	2.76	1.60	1.54
4	A	270	22B	C48-C47	2.77	1.57	1.52
3	E	260	RET	C18-C5	2.83	1.55	1.51
3	D	260	RET	C2-C1	2.86	1.60	1.54
4	B	270	22B	C48-C47	2.86	1.57	1.52
4	E	270	22B	C48-C47	2.87	1.57	1.52
4	D	270	22B	C48-C47	2.91	1.57	1.52
4	E	270	22B	C21-C2	3.14	1.57	1.53
4	D	270	22B	C34-C33	3.15	1.57	1.53
4	B	270	22B	C17-C1	3.18	1.58	1.52
4	A	270	22B	C17-C1	3.19	1.58	1.52
3	A	260	RET	C18-C5	3.19	1.56	1.51
4	D	270	22B	C17-C1	3.27	1.58	1.52
4	E	270	22B	C34-C33	3.35	1.58	1.53
4	E	270	22B	C17-C1	3.47	1.58	1.52
4	B	270	22B	C34-C33	3.50	1.58	1.53
4	A	270	22B	C34-C33	3.60	1.58	1.53
4	E	270	22B	O27-C23	3.64	1.54	1.44
4	B	270	22B	C16-C1	3.65	1.59	1.52
4	A	270	22B	C16-C1	3.70	1.59	1.52
4	E	270	22B	C49-C47	3.71	1.59	1.52
4	D	270	22B	C16-C1	3.75	1.59	1.52
4	D	270	22B	O27-C23	3.76	1.54	1.44
4	B	270	22B	O27-C23	3.77	1.54	1.44
4	A	270	22B	C49-C47	3.77	1.59	1.52
4	D	270	22B	C49-C47	3.79	1.59	1.52
4	B	270	22B	C49-C47	3.79	1.59	1.52
4	A	270	22B	O27-C23	3.80	1.54	1.44
3	A	260	RET	C7-C6	3.83	1.59	1.45
4	E	270	22B	C16-C1	4.10	1.59	1.52
3	B	260	RET	C7-C6	4.14	1.60	1.45
3	D	260	RET	C7-C6	4.19	1.60	1.45
3	E	260	RET	C7-C6	4.38	1.61	1.45
3	A	260	RET	C1-C6	4.58	1.60	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	260	RET	C1-C6	4.70	1.60	1.53
3	E	260	RET	C1-C6	5.09	1.60	1.53
3	B	260	RET	C1-C6	5.26	1.61	1.53
4	A	270	22B	C2-C3	6.64	1.57	1.49
4	B	270	22B	C2-C3	6.69	1.57	1.49
4	D	270	22B	C2-C3	7.16	1.57	1.49
4	B	270	22B	C33-C32	8.12	1.58	1.49
4	E	270	22B	C2-C3	8.24	1.59	1.49
4	D	270	22B	C33-C32	8.63	1.59	1.49
4	A	270	22B	C33-C32	8.72	1.59	1.49
3	A	260	RET	C5-C6	8.83	1.49	1.34
3	D	260	RET	C5-C6	8.87	1.49	1.34
4	E	270	22B	C33-C32	8.97	1.59	1.49
3	B	260	RET	C5-C6	9.17	1.50	1.34
3	E	260	RET	C5-C6	9.31	1.50	1.34

All (112) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	270	22B	C31-C29-C28	-4.68	111.75	118.94
4	E	270	22B	C31-C29-C28	-4.38	112.22	118.94
3	A	260	RET	C8-C9-C10	-3.82	113.08	118.94
3	D	260	RET	C8-C9-C10	-3.66	113.33	118.94
4	B	270	22B	C31-C29-C28	-3.59	113.43	118.94
3	B	260	RET	C8-C9-C10	-3.40	113.72	118.94
4	A	270	22B	C8-C9-C10	-3.30	113.88	118.94
4	B	270	22B	C8-C9-C10	-3.30	113.88	118.94
4	E	270	22B	C8-C9-C10	-3.29	113.90	118.94
4	D	270	22B	C8-C9-C10	-3.28	113.90	118.94
4	B	270	22B	C42-C43-C44	-3.16	114.09	118.94
4	A	270	22B	C42-C43-C44	-2.92	114.45	118.94
4	B	270	22B	C41-C40-C39	-2.91	123.16	127.31
4	D	270	22B	C41-C40-C39	-2.86	123.23	127.31
3	E	260	RET	C8-C9-C10	-2.84	114.58	118.94
4	E	270	22B	C42-C43-C44	-2.81	114.63	118.94
4	B	270	22B	C49-C47-C48	-2.76	106.42	110.53
4	A	270	22B	C31-C29-C28	-2.68	114.83	118.94
3	E	260	RET	C1-C6-C5	-2.68	118.83	122.59
4	A	270	22B	C17-C1-C16	-2.64	106.61	110.53
4	B	270	22B	C4-C5-C6	-2.64	114.90	118.94
3	D	260	RET	C18-C5-C6	-2.62	121.58	124.51
4	A	270	22B	C49-C47-C48	-2.61	106.65	110.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	270	22B	C42-C43-C44	-2.58	114.98	118.94
4	D	270	22B	C49-C47-C48	-2.57	106.70	110.53
4	D	270	22B	C12-C13-C14	-2.56	115.02	118.94
3	B	260	RET	C1-C6-C5	-2.55	119.00	122.59
4	E	270	22B	C49-C47-C48	-2.52	106.78	110.53
4	D	270	22B	C4-C5-C6	-2.52	115.08	118.94
3	D	260	RET	C1-C6-C5	-2.51	119.06	122.59
4	A	270	22B	C4-C5-C6	-2.48	115.13	118.94
3	E	260	RET	C18-C5-C6	-2.47	121.74	124.51
4	E	270	22B	C4-C5-C6	-2.47	115.15	118.94
4	D	270	22B	C17-C1-C16	-2.40	106.97	110.53
3	E	260	RET	C17-C1-C2	-2.38	99.43	108.80
3	A	260	RET	C1-C6-C5	-2.37	119.26	122.59
4	B	270	22B	C33-C32-C31	-2.34	121.25	125.18
4	B	270	22B	C17-C1-C16	-2.33	107.06	110.53
3	D	260	RET	C17-C1-C2	-2.26	99.87	108.80
3	A	260	RET	C17-C1-C2	-2.26	99.90	108.80
3	B	260	RET	C17-C1-C2	-2.24	99.98	108.80
3	A	260	RET	C18-C5-C6	-2.23	122.01	124.51
4	B	270	22B	C11-C12-C13	-2.21	120.21	126.42
3	B	260	RET	C10-C11-C12	-2.20	116.49	123.23
4	E	270	22B	C17-C1-C16	-2.20	107.26	110.53
3	D	260	RET	C10-C11-C12	-2.16	116.61	123.23
3	B	260	RET	C16-C1-C2	-2.14	100.37	108.80
4	E	270	22B	C12-C13-C14	-2.13	115.67	118.94
3	E	260	RET	C16-C1-C2	-2.12	100.44	108.80
3	D	260	RET	C16-C1-C2	-2.12	100.45	108.80
4	E	270	22B	C41-C40-C39	-2.12	124.29	127.31
3	A	260	RET	C16-C1-C2	-2.10	100.50	108.80
4	A	270	22B	C41-C40-C39	-2.10	124.32	127.31
4	A	270	22B	C11-C12-C13	-2.06	120.63	126.42
3	A	260	RET	C10-C11-C12	-2.03	117.02	123.23
3	B	260	RET	C18-C5-C6	-2.02	122.24	124.51
4	B	270	22B	C12-C13-C14	-2.02	115.84	118.94
4	D	270	22B	C26-C39-C38	2.15	121.53	118.10
3	E	260	RET	C3-C2-C1	2.16	122.38	114.59
4	A	270	22B	C26-C39-C38	2.17	121.56	118.10
4	E	270	22B	C20-C13-C12	2.20	121.61	118.10
4	D	270	22B	C20-C13-C12	2.22	121.63	118.10
3	A	260	RET	C3-C2-C1	2.22	122.58	114.59
3	B	260	RET	C20-C13-C12	2.26	121.70	118.10
4	E	270	22B	C26-C39-C38	2.35	121.84	118.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	260	RET	C20-C13-C12	2.35	121.84	118.10
3	A	260	RET	C20-C13-C12	2.37	121.87	118.10
3	E	260	RET	C19-C9-C8	2.37	121.88	118.10
3	E	260	RET	C20-C13-C12	2.45	122.00	118.10
3	D	260	RET	C19-C9-C8	2.47	122.03	118.10
3	B	260	RET	C2-C3-C4	2.51	117.32	111.34
4	D	270	22B	C50-C43-C42	2.51	122.10	118.10
3	A	260	RET	C1-C6-C7	2.53	122.83	115.73
3	E	260	RET	C17-C1-C6	2.53	114.40	110.31
3	B	260	RET	C1-C6-C7	2.55	122.90	115.73
3	D	260	RET	C2-C3-C4	2.55	117.43	111.34
4	A	270	22B	C30-C29-C31	2.56	122.17	118.10
4	E	270	22B	C50-C43-C42	2.57	122.20	118.10
3	A	260	RET	C19-C9-C8	2.60	122.24	118.10
3	B	260	RET	C19-C9-C8	2.62	122.27	118.10
4	A	270	22B	C50-C43-C42	2.68	122.37	118.10
4	E	270	22B	C18-C5-C4	2.72	122.43	118.10
4	B	270	22B	C50-C43-C42	2.74	122.47	118.10
3	D	260	RET	C1-C6-C7	2.74	123.44	115.73
4	B	270	22B	C18-C5-C4	2.74	122.47	118.10
4	D	270	22B	C18-C5-C4	2.76	122.49	118.10
3	E	260	RET	C1-C6-C7	2.77	123.51	115.73
3	D	260	RET	C17-C1-C6	2.81	114.86	110.31
4	A	270	22B	C18-C5-C4	2.81	122.57	118.10
3	B	260	RET	C17-C1-C6	2.84	114.92	110.31
4	E	270	22B	C19-C9-C8	2.85	122.64	118.10
3	B	260	RET	C3-C4-C5	2.86	118.69	113.78
3	A	260	RET	C17-C1-C6	2.91	115.02	110.31
3	D	260	RET	C3-C4-C5	2.92	118.81	113.78
4	A	270	22B	C19-C9-C8	2.95	122.80	118.10
4	D	270	22B	C19-C9-C8	2.96	122.81	118.10
3	A	260	RET	C16-C1-C6	2.96	115.11	110.31
4	B	270	22B	C30-C29-C31	2.98	122.85	118.10
3	A	260	RET	C2-C3-C4	3.01	118.52	111.34
4	B	270	22B	C19-C9-C8	3.02	122.91	118.10
3	E	260	RET	C2-C3-C4	3.02	118.54	111.34
3	B	260	RET	C16-C1-C6	3.07	115.29	110.31
3	D	260	RET	C16-C1-C6	3.10	115.33	110.31
3	E	260	RET	C16-C1-C6	3.18	115.47	110.31
3	D	260	RET	C2-C1-C6	3.22	115.51	110.48
3	B	260	RET	C2-C1-C6	3.32	115.66	110.48
3	E	260	RET	C3-C4-C5	3.34	119.53	113.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	270	22B	C30-C29-C31	3.41	123.53	118.10
3	A	260	RET	C2-C1-C6	3.44	115.86	110.48
4	D	270	22B	C30-C29-C31	3.55	123.75	118.10
3	A	260	RET	C3-C4-C5	3.56	119.90	113.78
3	E	260	RET	C2-C1-C6	3.70	116.26	110.48

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

10 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	260	RET	2	0
4	A	270	22B	3	0
5	A	280	L2P	5	0
3	B	260	RET	1	0
4	B	270	22B	2	0
5	B	280	L2P	1	0
4	D	270	22B	2	0
5	D	280	L2P	3	0
3	E	260	RET	2	0
5	E	280	L2P	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	242/253 (95%)	-0.38	1 (0%) 92 92	20, 29, 56, 78	0
1	B	238/253 (94%)	-0.38	1 (0%) 92 92	22, 34, 53, 68	0
1	D	237/253 (93%)	-0.33	4 (1%) 70 72	26, 43, 63, 93	0
1	E	238/253 (94%)	-0.23	5 (2%) 64 66	31, 51, 73, 85	0
All	All	955/1012 (94%)	-0.33	11 (1%) 79 80	20, 39, 67, 93	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	240	ALA	4.4
1	D	240	ALA	3.9
1	D	238	THR	3.4
1	D	241	PRO	3.2
1	E	238	THR	2.7
1	B	3	GLY	2.6
1	D	80	VAL	2.3
1	E	237	GLU	2.1
1	E	228	LEU	2.1
1	E	231	SER	2.1
1	E	227	VAL	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](#)

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	MAN	B	282	11/12	0.82	0.44	11.81	106,108,109,109	0
2	GAL	A	283	11/12	0.88	0.38	7.04	112,112,113,114	0
2	MAN	A	282	11/12	0.60	0.29	4.48	109,109,110,111	0
2	GAL	D	283	11/12	0.79	0.28	4.14	108,109,109,109	0
2	MAN	D	282	11/12	0.61	0.23	3.05	106,107,108,108	0
2	GAL	B	283	11/12	0.82	0.35	2.27	109,110,110,111	0
2	MAN	D	285	11/12	0.70	0.31	1.85	108,110,110,110	0
2	GAL	D	286	11/12	0.84	0.28	0.81	109,109,110,110	0
2	GLC	A	281	11/12	0.62	0.26	-	104,106,107,107	0
2	GLC	B	281	11/12	0.85	0.19	-	102,102,103,103	0
2	GLC	D	281	11/12	0.83	0.10	-	104,104,105,105	0
2	GLC	D	284	11/12	0.72	0.17	-	106,107,107,107	0

## 6.4 Ligands ⓘ

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
5	L2P	E	280	46/46	0.61	0.53	19.04	95,99,102,104	0
5	L2P	A	280	46/46	0.69	0.50	16.38	84,98,104,105	0
5	L2P	D	280	46/46	0.65	0.60	16.29	89,101,103,103	0
5	L2P	B	280	46/46	0.67	0.50	13.88	80,98,100,101	0
4	22B	E	270	54/54	0.74	0.54	11.66	89,94,101,102	0
4	22B	D	270	54/54	0.78	0.41	9.28	79,87,103,104	0
4	22B	B	270	54/54	0.85	0.33	8.92	57,75,93,94	0
4	22B	A	270	54/54	0.87	0.27	6.46	63,72,94,95	0
3	RET	E	260	20/21	0.89	0.24	3.91	58,59,62,62	0
3	RET	D	260	20/21	0.94	0.19	3.17	30,34,36,38	0
3	RET	B	260	20/21	0.96	0.18	1.59	26,30,33,33	0
3	RET	A	260	20/21	0.94	0.15	1.32	20,22,27,27	0



## 6.5 Other polymers [i](#)

There are no such residues in this entry.