



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:10 AM GMT

PDB ID : 1ZUM  
Title : Human Mitochondrial Aldehyde Dehydrogenase Asian Variant, ALDH2\*2, Apo Form  
Authors : Larson, H.N.; Weiner, H.; Hurley, T.D.  
Deposited on : 2005-05-31  
Resolution : 2.10 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 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

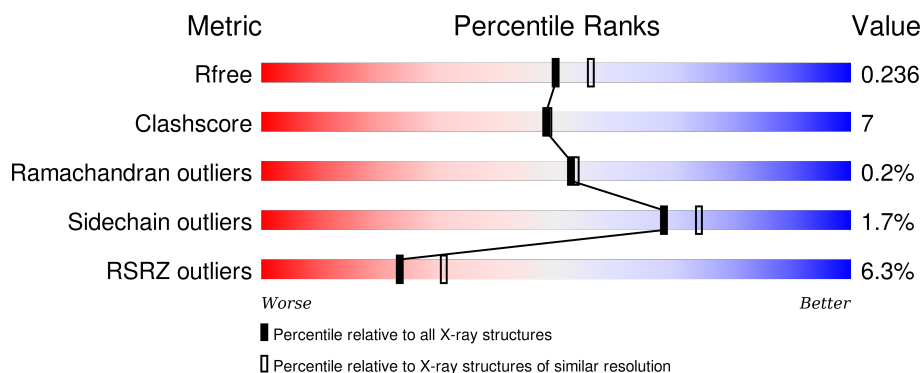
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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))
$R_{free}$	91344	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	500	<div> <div>8%</div> <div> <div></div> <div>80%</div> <div>13%</div> <div>6%</div> </div> </div>
1	B	500	<div> <div>2%</div> <div> <div></div> <div>83%</div> <div>8%</div> <div>7%</div> </div> </div>
1	C	500	<div> <div>2%</div> <div> <div></div> <div>80%</div> <div>12%</div> <div>6%</div> </div> </div>
1	D	500	<div> <div>8%</div> <div> <div></div> <div>76%</div> <div>17%</div> <div>6%</div> </div> </div>
1	E	500	<div> <div>3%</div> <div> <div></div> <div>82%</div> <div>11%</div> <div>6%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	500	
1	G	500	
1	H	500	
1	I	500	
1	J	500	
1	K	500	
1	L	500	

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
3	GAI	A	3101	-	-	X	-
3	GAI	F	3103	-	-	-	X
3	GAI	I	3107	-	-	X	X
3	GAI	L	3108	-	-	X	X
4	EDO	A	3109	-	-	-	X
4	EDO	A	3110	-	-	-	X
4	EDO	A	3116	-	-	X	X
4	EDO	C	3113	-	-	-	X
4	EDO	C	3114	-	-	-	X
4	EDO	D	3126	-	-	X	-
4	EDO	F	3120	-	-	-	X
4	EDO	F	3122	-	-	-	X
4	EDO	G	3123	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 46002 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 Aldehyde dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	469	Total	C	N	O	S	0	0	0
			3620	2306	618	679	17			
1	B	465	Total	C	N	O	S	0	0	0
			3585	2286	610	672	17			
1	C	468	Total	C	N	O	S	0	0	0
			3612	2302	617	676	17			
1	D	468	Total	C	N	O	S	0	0	0
			3608	2301	613	677	17			
1	E	469	Total	C	N	O	S	0	0	0
			3624	2311	618	678	17			
1	F	468	Total	C	N	O	S	0	0	0
			3612	2302	617	676	17			
1	G	469	Total	C	N	O	S	0	0	0
			3624	2311	618	678	17			
1	H	469	Total	C	N	O	S	0	0	0
			3624	2311	618	678	17			
1	I	468	Total	C	N	O	S	0	0	0
			3615	2305	616	677	17			
1	J	467	Total	C	N	O	S	0	0	0
			3604	2299	612	676	17			
1	K	456	Total	C	N	O	S	0	0	0
			3526	2250	598	661	17			
1	L	462	Total	C	N	O	S	0	0	0
			3571	2279	607	668	17			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	487	LYS	GLU	ENGINEERED	UNP P05091
B	487	LYS	GLU	ENGINEERED	UNP P05091
C	487	LYS	GLU	ENGINEERED	UNP P05091
D	487	LYS	GLU	ENGINEERED	UNP P05091
E	487	LYS	GLU	ENGINEERED	UNP P05091

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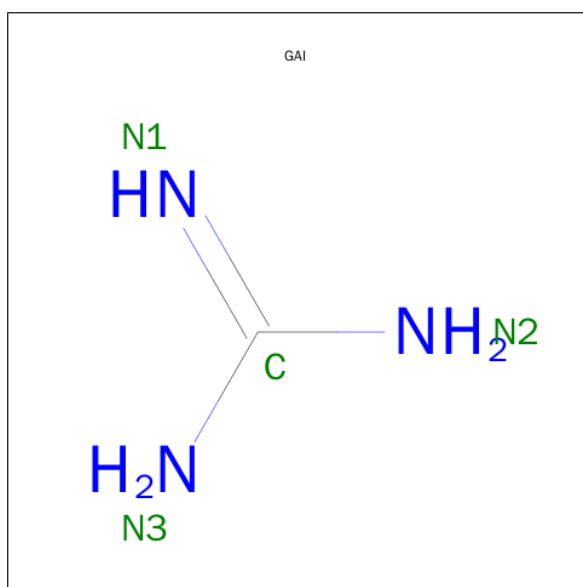
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Chain	Residue	Modelled	Actual	Comment	Reference
F	487	LYS	GLU	ENGINEERED	UNP P05091
G	487	LYS	GLU	ENGINEERED	UNP P05091
H	487	LYS	GLU	ENGINEERED	UNP P05091
I	487	LYS	GLU	ENGINEERED	UNP P05091
J	487	LYS	GLU	ENGINEERED	UNP P05091
K	487	LYS	GLU	ENGINEERED	UNP P05091
L	487	LYS	GLU	ENGINEERED	UNP P05091

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	G	1	Total Na 1 1	0	0
2	J	1	Total Na 1 1	0	0
2	D	1	Total Na 1 1	0	0
2	K	1	Total Na 1 1	0	0
2	E	1	Total Na 1 1	0	0
2	H	1	Total Na 1 1	0	0
2	B	1	Total Na 1 1	0	0
2	I	1	Total Na 1 1	0	0
2	C	1	Total Na 1 1	0	0
2	A	1	Total Na 1 1	0	0
2	L	1	Total Na 1 1	0	0
2	F	1	Total Na 1 1	0	0

- Molecule 3 is GUANIDINE (three-letter code: GAI) (formula: CH<sub>5</sub>N<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			4	1	3		
3	E	1	Total	C	N	0	0
			4	1	3		
3	F	1	Total	C	N	0	0
			4	1	3		
3	F	1	Total	C	N	0	0
			4	1	3		
3	H	1	Total	C	N	0	0
			4	1	3		
3	I	1	Total	C	N	0	0
			4	1	3		
3	L	1	Total	C	N	0	0
			4	1	3		

- Molecule 4 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
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		
4	D	1	Total	C	O	0	0
			4	2	2		
4	E	1	Total	C	O	0	0
			4	2	2		
4	I	1	Total	C	O	0	0
			4	2	2		
4	F	1	Total	C	O	0	0
			4	2	2		
4	F	1	Total	C	O	0	0
			4	2	2		
4	F	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	G	1	Total C O 4 2 2	0	0
4	G	1	Total C O 4 2 2	0	0
4	G	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	I	1	Total C O 4 2 2	0	0
4	J	1	Total C O 4 2 2	0	0
4	L	1	Total C O 4 2 2	0	0

- Molecule 5 is water.

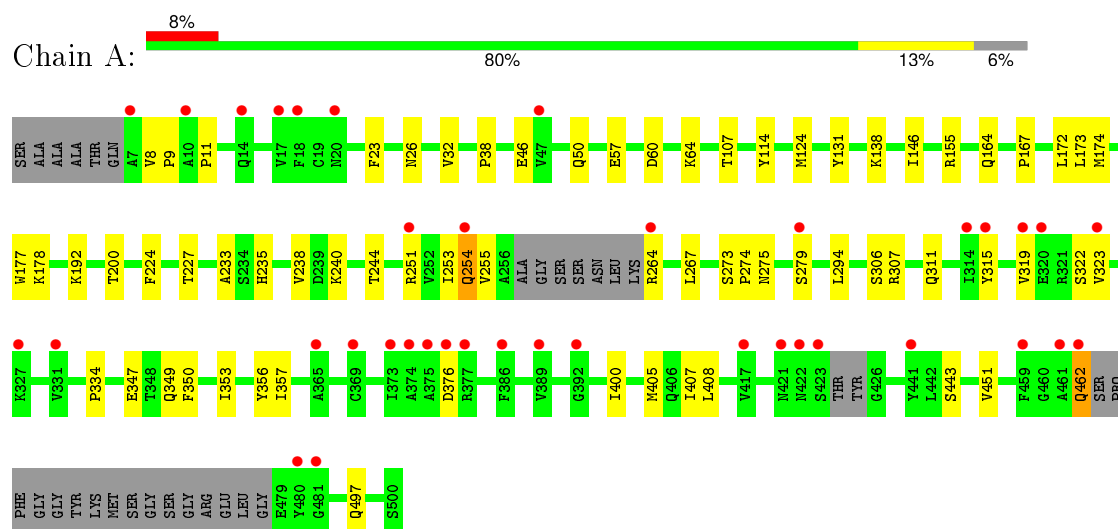
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	189	Total O 189 189	0	0
5	B	299	Total O 299 299	0	0
5	C	290	Total O 290 290	0	0
5	D	187	Total O 187 187	0	0
5	E	285	Total O 285 285	0	0
5	F	367	Total O 367 367	0	0
5	G	285	Total O 285 285	0	0
5	H	236	Total O 236 236	0	0
5	I	189	Total O 189 189	0	0
5	J	141	Total O 141 141	0	0
5	K	93	Total O 93 93	0	0
5	L	88	Total O 88 88	0	0



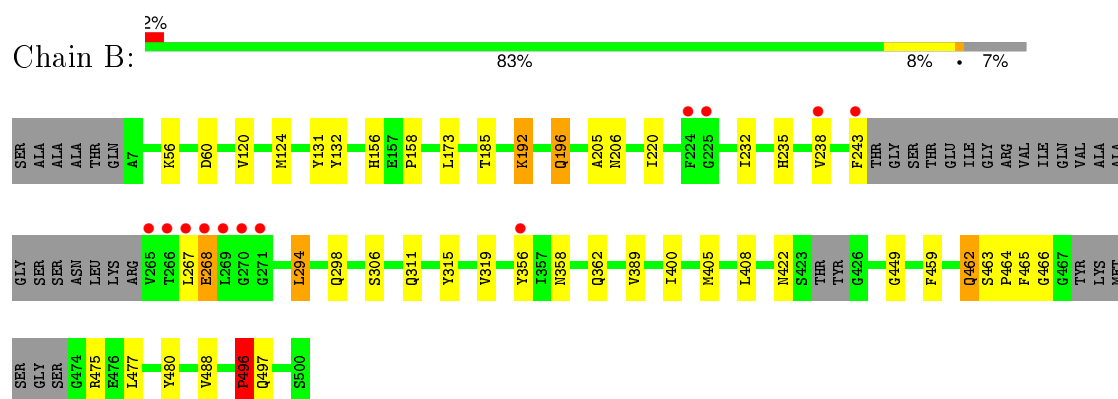
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

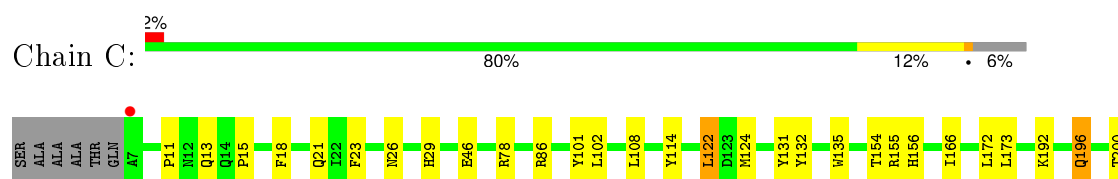
#### • Molecule 1: Aldehyde dehydrogenase

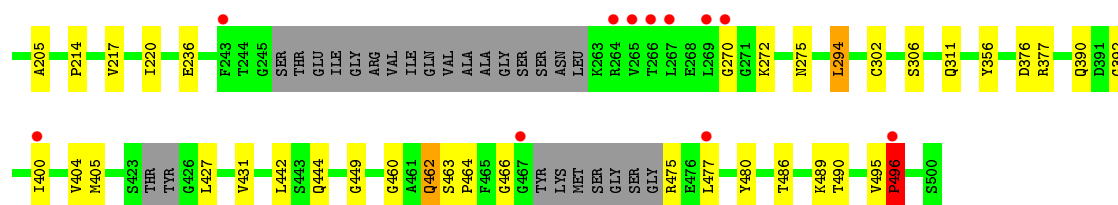


#### • Molecule 1: Aldehyde dehydrogenase

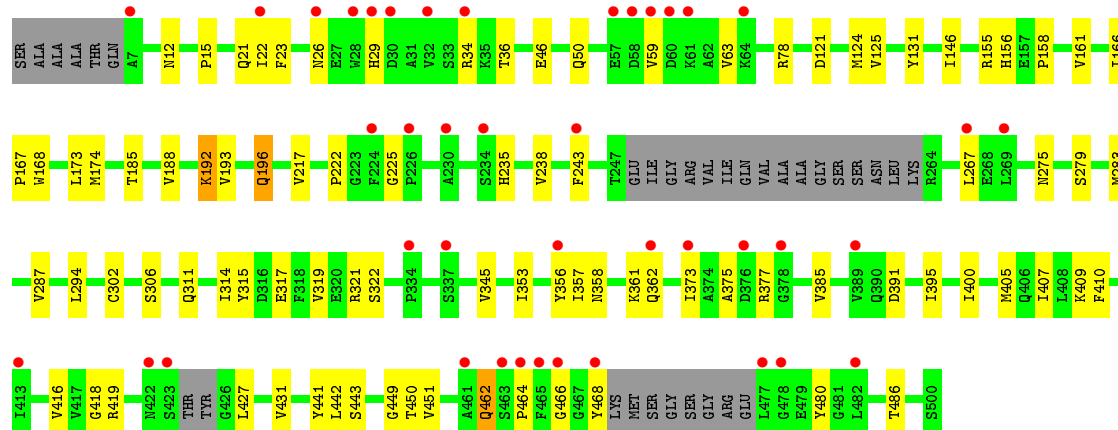
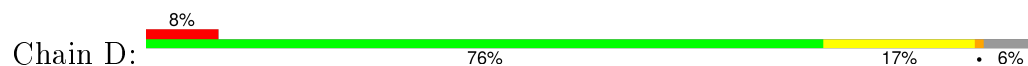


#### • Molecule 1: Aldehyde dehydrogenase

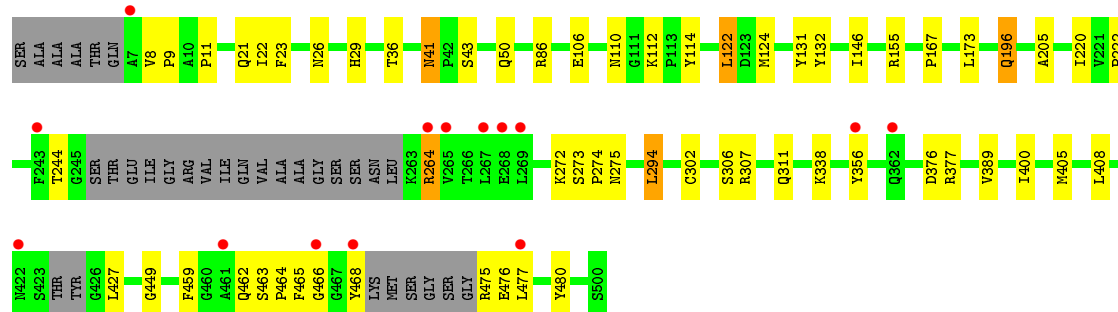
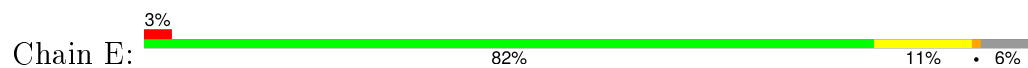




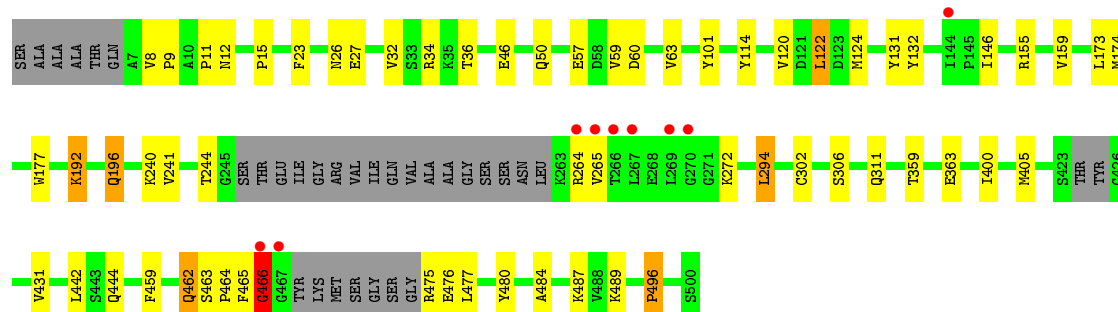
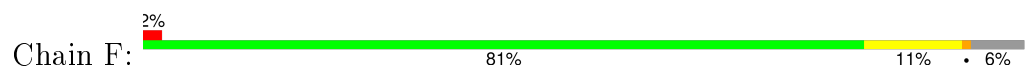
• Molecule 1: Aldehyde dehydrogenase



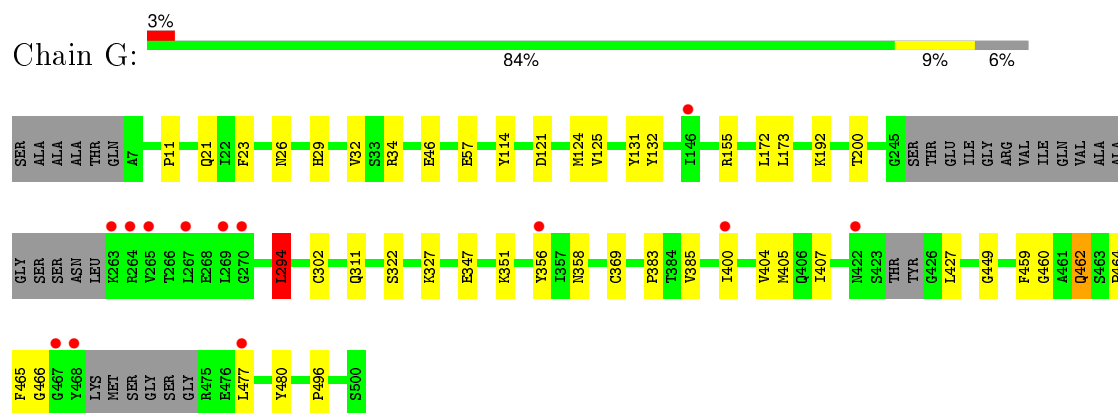
• Molecule 1: Aldehyde dehydrogenase



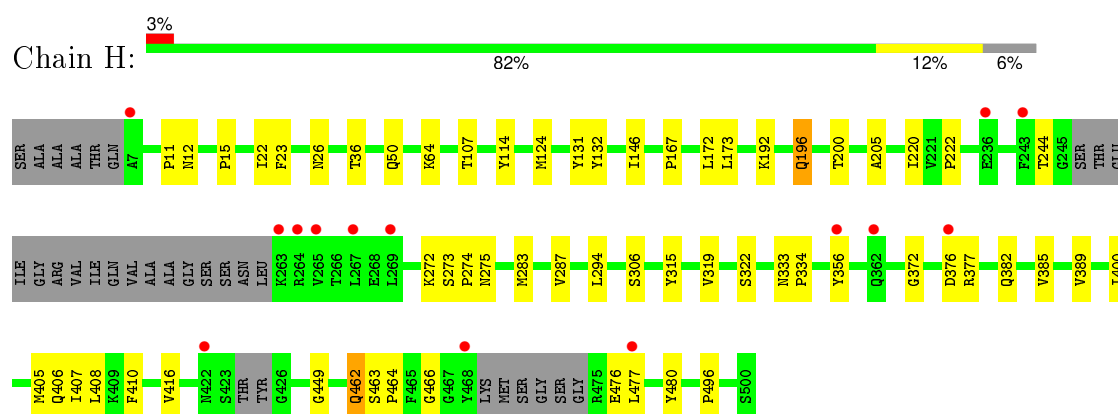
• Molecule 1: Aldehyde dehydrogenase



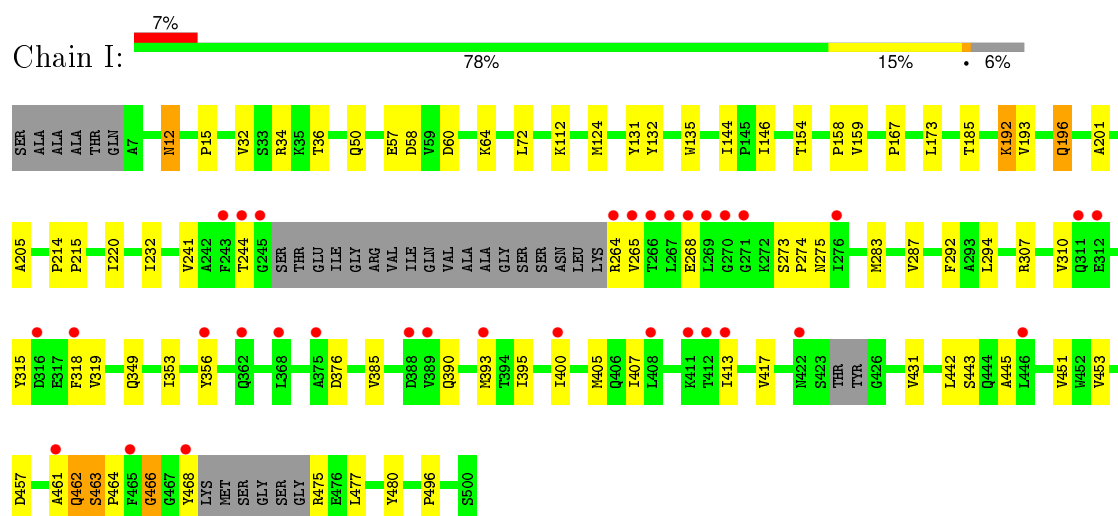
- Molecule 1: Aldehyde dehydrogenase



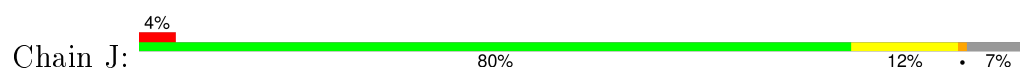
- Molecule 1: Aldehyde dehydrogenase

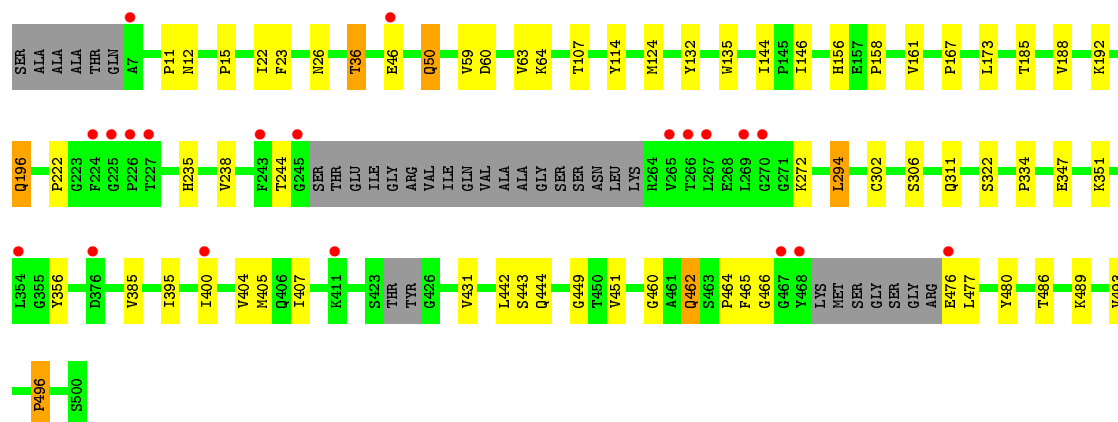


- Molecule 1: Aldehyde dehydrogenase

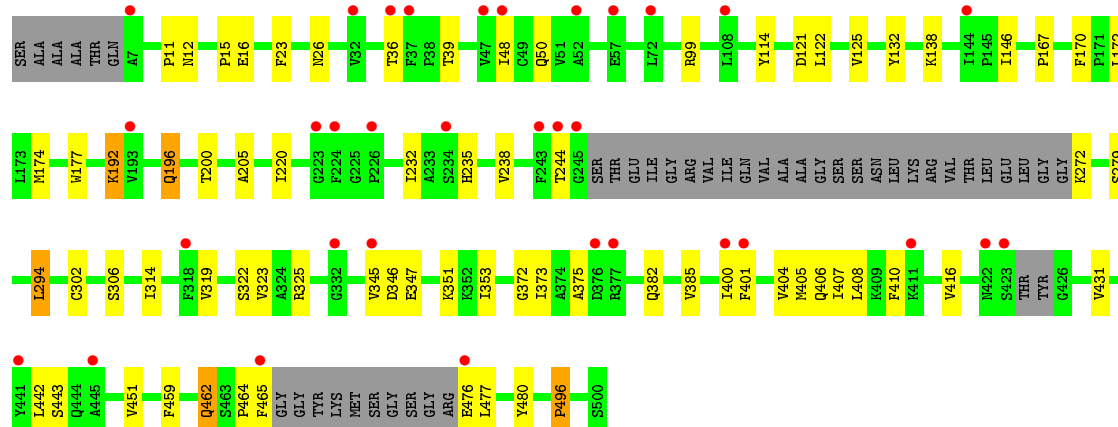
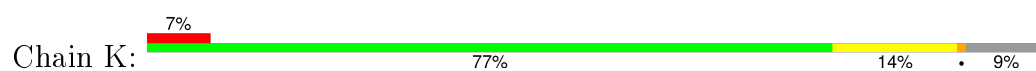


- Molecule 1: Aldehyde dehydrogenase

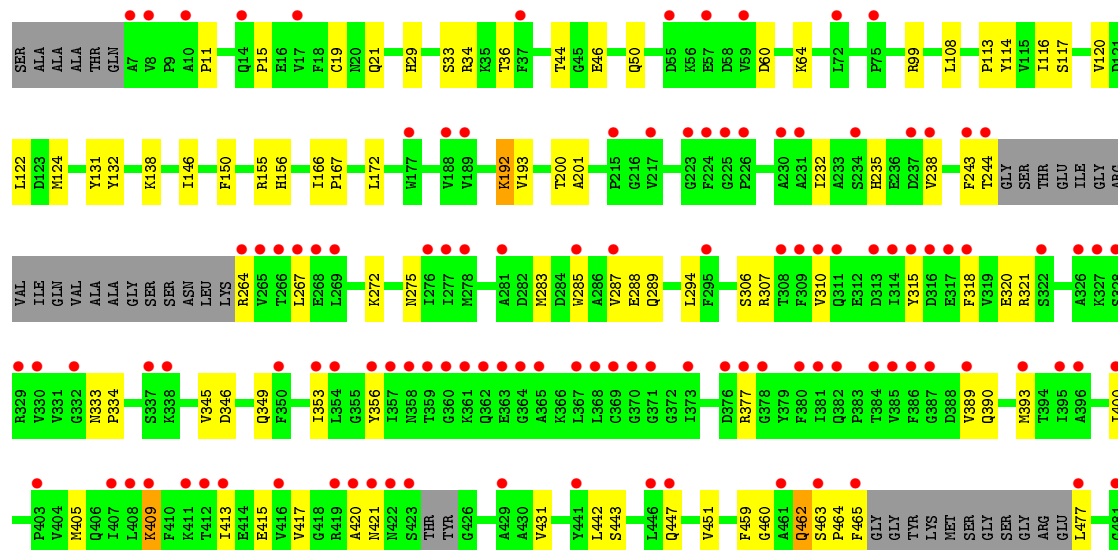


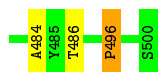


• Molecule 1: Aldehyde dehydrogenase



• Molecule 1: Aldehyde dehydrogenase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	95.87Å 105.26Å 162.67Å 78.81° 81.95° 88.04°	Depositor
Resolution (Å)	42.63 – 2.10 42.62 – 2.11	Depositor EDS
% Data completeness (in resolution range)	97.7 (42.63-2.10) 91.5 (42.62-2.11)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.33 (at 2.10Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.204 , 0.238 0.205 , 0.236	Depositor DCC
$R_{free}$ test set	17348 reflections (5.03%)	DCC
Wilson B-factor (Å <sup>2</sup> )	31.4	Xtriage
Anisotropy	0.266	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 49.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 345156 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	46002	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.84% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: NA, EDO, GAI

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	0.41	0/3697	0.61	0/5015
1	B	0.51	0/3664	0.67	0/4969
1	C	0.52	0/3691	0.69	2/5004 (0.0%)
1	D	0.42	0/3688	0.61	0/5003
1	E	0.51	0/3704	0.68	1/5022 (0.0%)
1	F	0.56	0/3691	0.69	3/5004 (0.1%)
1	G	0.50	0/3704	0.66	1/5022 (0.0%)
1	H	0.50	0/3704	0.65	0/5022
1	I	0.41	0/3695	0.62	1/5011 (0.0%)
1	J	0.38	0/3684	0.60	0/4997
1	K	0.36	0/3605	0.58	0/4891
1	L	0.36	0/3650	0.58	0/4952
All	All	0.46	0/44177	0.64	8/59912 (0.0%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	122	LEU	CA-CB-CG	-7.76	97.44	115.30
1	C	122	LEU	CA-CB-CG	-7.38	98.31	115.30
1	F	122	LEU	CA-CB-CG	-6.86	99.53	115.30
1	C	294	LEU	CA-CB-CG	-6.34	100.71	115.30
1	G	294	LEU	CA-CB-CG	-5.67	102.25	115.30
1	F	466	GLY	N-CA-C	5.44	126.70	113.10
1	F	294	LEU	CA-CB-CG	-5.35	103.00	115.30
1	I	12	ASN	N-CA-C	-5.08	97.28	111.00

There are no chirality outliers.

There are no planarity outliers.

## 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	3620	0	3578	53	0
1	B	3585	0	3531	42	0
1	C	3612	0	3564	57	0
1	D	3608	0	3553	72	0
1	E	3624	0	3573	59	0
1	F	3612	0	3564	57	0
1	G	3624	0	3573	38	0
1	H	3624	0	3573	46	0
1	I	3615	0	3560	65	0
1	J	3604	0	3547	51	0
1	K	3526	0	3469	51	0
1	L	3571	0	3523	64	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	K	1	0	0	0	0
2	L	1	0	0	0	0
3	A	4	0	4	2	0
3	E	4	0	5	0	0
3	F	8	0	10	1	0
3	G	4	0	4	0	0
3	H	4	0	5	0	0
3	I	4	0	4	2	0
3	L	4	0	5	4	0
4	A	16	0	24	8	0
4	B	4	0	6	0	0
4	C	12	0	18	5	0
4	D	8	0	12	6	0
4	E	4	0	6	0	0
4	F	12	0	18	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	G	12	0	18	2	0
4	I	8	0	12	0	0
4	J	4	0	6	1	0
4	L	4	0	6	1	0
5	A	189	0	0	3	0
5	B	299	0	0	5	0
5	C	290	0	0	5	0
5	D	187	0	0	5	0
5	E	285	0	0	6	0
5	F	367	0	0	9	0
5	G	285	0	0	6	0
5	H	236	0	0	2	0
5	I	189	0	0	4	0
5	J	141	0	0	3	0
5	K	93	0	0	1	0
5	L	88	0	0	6	0
All	All	46002	0	42771	629	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 7.

All (629) 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:H:124:MET:HE3	1:H:173:LEU:HD22	1.52	0.89
1:C:196:GLN:HE21	1:C:196:GLN:H	1.20	0.88
1:I:466:GLY:HA3	1:I:475:ARG:HD3	1.53	0.87
1:A:131:TYR:CE1	1:A:462:GLN:HG3	2.11	0.86
1:L:36:THR:OG1	1:L:50:GLN:HG3	1.75	0.86
1:H:462:GLN:H	1:H:462:GLN:NE2	1.75	0.85
1:L:421:ASN:HD21	1:L:447:GLN:HB2	1.44	0.83
1:I:196:GLN:HE21	1:I:196:GLN:H	1.29	0.80
1:D:462:GLN:H	1:D:462:GLN:NE2	1.79	0.80
1:K:196:GLN:H	1:K:196:GLN:HE21	1.28	0.79
1:G:124:MET:HE3	1:G:173:LEU:HD22	1.64	0.79
1:L:390:GLN:H	1:L:393:MET:HE2	1.44	0.79
1:E:196:GLN:H	1:E:196:GLN:HE21	1.29	0.79
1:F:196:GLN:H	1:F:196:GLN:HE21	1.31	0.79
1:F:155:ARG:HD2	4:F:3120:EDO:O2	1.83	0.79
1:H:196:GLN:HE21	1:H:196:GLN:H	1.31	0.78
1:F:462:GLN:NE2	1:F:462:GLN:H	1.82	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:294:LEU:HD12	1:L:306:SER:HA	1.64	0.78
1:F:302:CYS:HB3	5:F:4036:HOH:O	1.84	0.77
1:G:302:CYS:HB3	5:G:3968:HOH:O	1.85	0.77
1:I:124:MET:HE3	1:I:173:LEU:HD22	1.65	0.76
1:J:124:MET:HE3	1:J:173:LEU:HD22	1.67	0.76
1:E:124:MET:HE3	1:E:173:LEU:HD22	1.67	0.76
1:G:462:GLN:NE2	1:G:462:GLN:H	1.84	0.76
1:B:196:GLN:H	1:B:196:GLN:HE21	1.32	0.76
4:A:3116:EDO:H11	1:C:444:GLN:CG	2.17	0.75
1:C:302:CYS:HB3	5:C:3956:HOH:O	1.86	0.75
1:K:36:THR:OG1	1:K:50:GLN:HG3	1.86	0.74
1:A:279:SER:H	1:A:311:GLN:HE21	1.35	0.74
1:L:155:ARG:HD2	4:L:3129:EDO:O2	1.86	0.74
1:I:159:VAL:HG21	1:I:264:ARG:NH1	2.03	0.73
1:D:124:MET:HE3	1:D:173:LEU:HD22	1.70	0.73
1:L:421:ASN:ND2	1:L:447:GLN:HB2	2.03	0.73
1:A:279:SER:H	1:A:311:GLN:NE2	1.86	0.73
1:F:124:MET:HE3	1:F:173:LEU:HD22	1.70	0.73
1:E:338:LYS:HD2	1:I:34:ARG:HH11	1.53	0.73
1:K:205:ALA:HB2	1:K:220:ILE:HD12	1.69	0.73
1:G:400:ILE:HD11	1:G:404:VAL:HB	1.68	0.72
1:E:41:ASN:ND2	1:E:43:SER:H	1.88	0.71
1:F:466:GLY:HA3	1:F:475:ARG:HD3	1.72	0.71
1:D:196:GLN:H	1:D:196:GLN:HE21	1.36	0.70
1:F:487:LYS:HD3	5:F:3915:HOH:O	1.91	0.70
1:F:60:ASP:HB2	5:F:4019:HOH:O	1.90	0.70
1:B:497:GLN:HG3	1:C:78:ARG:NH1	2.07	0.70
1:A:294:LEU:HD12	1:A:306:SER:HA	1.72	0.70
1:H:294:LEU:HD12	1:H:306:SER:HA	1.74	0.69
1:B:462:GLN:NE2	1:B:462:GLN:H	1.90	0.69
1:F:294:LEU:HD22	1:F:405:MET:HB2	1.72	0.69
1:E:264:ARG:HH11	1:E:264:ARG:HB2	1.58	0.69
1:B:311:GLN:HG2	5:B:3721:HOH:O	1.91	0.69
1:J:36:THR:OG1	1:J:50:GLN:HG3	1.92	0.69
1:D:36:THR:OG1	1:D:50:GLN:HG3	1.93	0.69
1:J:196:GLN:H	1:J:196:GLN:HE21	1.41	0.69
1:L:390:GLN:H	1:L:393:MET:CE	2.06	0.69
1:J:302:CYS:HB3	5:J:3829:HOH:O	1.92	0.68
1:D:311:GLN:HG2	5:D:3753:HOH:O	1.93	0.68
1:B:464:PRO:HG3	1:B:480:TYR:CD1	2.27	0.68
1:F:240:LYS:HG3	1:F:264:ARG:HD2	1.74	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:462:GLN:H	1:L:462:GLN:NE2	1.92	0.68
1:A:240:LYS:HE2	1:A:264:ARG:NH1	2.08	0.68
1:J:462:GLN:H	1:J:462:GLN:NE2	1.93	0.67
1:C:196:GLN:NE2	1:C:196:GLN:H	1.90	0.67
1:L:389:VAL:HA	1:L:393:MET:HE1	1.77	0.66
1:K:170:PHE:HB2	1:K:174:MET:HE3	1.78	0.66
1:G:155:ARG:HD2	4:G:3123:EDO:O2	1.95	0.66
1:K:146:ILE:HG12	3:L:3108:GAI:HN31	1.60	0.66
1:E:462:GLN:HG2	1:F:146:ILE:HG22	1.78	0.66
1:I:468:TYR:OH	1:J:489:LYS:HE3	1.96	0.66
4:A:3116:EDO:O2	1:D:155:ARG:HD2	1.95	0.66
1:J:311:GLN:HG2	5:J:3825:HOH:O	1.95	0.66
1:L:272:LYS:HG3	1:L:307:ARG:HD3	1.78	0.65
1:D:302:CYS:SG	5:D:3865:HOH:O	2.53	0.65
1:E:377:ARG:HG3	1:E:377:ARG:HH11	1.59	0.65
1:L:99:ARG:HG3	1:L:122:LEU:HD23	1.76	0.65
1:E:294:LEU:HD22	1:E:405:MET:HB2	1.77	0.65
1:D:294:LEU:HD12	1:D:306:SER:HA	1.77	0.65
1:D:302:CYS:HB3	5:D:3864:HOH:O	1.95	0.65
1:F:302:CYS:SG	5:F:4035:HOH:O	2.54	0.65
1:C:294:LEU:HD22	1:C:405:MET:HB2	1.80	0.64
1:E:41:ASN:C	1:E:41:ASN:HD22	2.00	0.64
1:F:462:GLN:NE2	5:F:3953:HOH:O	2.30	0.64
1:I:462:GLN:NE2	1:I:462:GLN:H	1.96	0.64
1:A:462:GLN:NE2	1:A:462:GLN:H	1.95	0.63
4:A:3116:EDO:H11	1:C:444:GLN:HG3	1.80	0.63
1:G:322:SER:HB3	1:G:405:MET:HE3	1.80	0.63
1:I:132:TYR:OH	1:I:477:LEU:HA	1.98	0.63
1:C:462:GLN:NE2	1:C:462:GLN:H	1.95	0.63
1:J:156:HIS:HB3	1:J:486:THR:HG21	1.80	0.63
1:C:311:GLN:HG2	5:C:3748:HOH:O	1.98	0.62
1:C:78:ARG:HG2	1:C:78:ARG:HH11	1.64	0.62
4:A:3116:EDO:H11	1:C:444:GLN:HG2	1.81	0.62
1:K:294:LEU:HD22	1:K:405:MET:HB2	1.80	0.62
1:E:311:GLN:HG2	5:E:3756:HOH:O	1.99	0.62
1:A:405:MET:HE2	1:A:407:ILE:HD11	1.81	0.62
1:K:99:ARG:HG3	1:K:122:LEU:HD23	1.82	0.61
1:D:405:MET:HE3	1:D:407:ILE:HD11	1.82	0.61
1:C:124:MET:HE3	1:C:173:LEU:HD22	1.82	0.61
1:H:196:GLN:H	1:H:196:GLN:NE2	1.99	0.61
1:D:168:TRP:CE3	1:D:345:VAL:HG21	2.36	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:GLU:HB2	4:A:3110:EDO:H21	1.83	0.61
1:F:196:GLN:H	1:F:196:GLN:NE2	1.98	0.61
1:D:431:VAL:HG21	1:D:442:LEU:HB3	1.83	0.60
1:B:235:HIS:HB3	1:B:238:VAL:HG23	1.83	0.60
1:A:60:ASP:O	1:A:64:LYS:HG2	2.01	0.60
1:C:156:HIS:HB3	1:C:486:THR:HG21	1.83	0.60
1:L:19:CYS:HA	5:L:3769:HOH:O	2.01	0.60
5:A:3864:HOH:O	1:C:86:ARG:HD3	2.00	0.60
1:E:464:PRO:HG3	1:E:480:TYR:CD1	2.36	0.60
1:K:462:GLN:HG2	1:L:146:ILE:HG22	1.83	0.60
1:E:338:LYS:HD2	1:I:34:ARG:NH1	2.16	0.60
1:H:22:ILE:HD13	1:H:222:PRO:HD2	1.83	0.60
1:L:272:LYS:HG3	1:L:307:ARG:CD	2.31	0.60
1:A:155:ARG:HD2	4:A:3109:EDO:O2	1.99	0.60
1:A:124:MET:HE3	1:A:173:LEU:HD22	1.82	0.60
1:D:12:ASN:O	1:D:15:PRO:HD3	2.01	0.60
1:I:159:VAL:HG21	1:I:264:ARG:HH12	1.66	0.59
1:F:36:THR:HB	1:F:50:GLN:HG3	1.83	0.59
1:J:294:LEU:HD12	1:J:306:SER:HA	1.84	0.59
1:K:177:TRP:CD1	1:K:477:LEU:HD21	2.37	0.59
1:D:317:GLU:HG2	1:D:321:ARG:HE	1.67	0.59
1:B:124:MET:HE3	1:B:173:LEU:HD22	1.83	0.59
1:E:356:TYR:CD2	1:E:400:ILE:HG12	2.38	0.59
1:D:294:LEU:HD13	1:D:405:MET:HA	1.85	0.59
1:F:464:PRO:HG3	1:F:480:TYR:HD1	1.68	0.59
1:G:356:TYR:CD1	1:G:400:ILE:HG22	2.38	0.58
1:E:41:ASN:HD22	1:E:43:SER:H	1.50	0.58
1:E:462:GLN:NE2	5:E:3929:HOH:O	2.36	0.58
1:C:460:GLY:HA3	1:D:146:ILE:HG13	1.84	0.58
1:J:60:ASP:O	1:J:64:LYS:HG2	2.02	0.58
1:D:385:VAL:HG13	1:D:405:MET:HE2	1.85	0.58
1:K:172:LEU:HD21	1:K:200:THR:HB	1.85	0.58
1:J:465:PHE:O	1:J:476:GLU:HB2	2.02	0.58
1:L:44:THR:OG1	1:L:46:GLU:HG2	2.04	0.58
1:J:196:GLN:H	1:J:196:GLN:NE2	2.01	0.58
1:I:376:ASP:HA	5:I:3746:HOH:O	2.04	0.58
1:L:172:LEU:HD21	1:L:200:THR:HB	1.86	0.57
1:B:120:VAL:HG12	1:B:124:MET:HE1	1.86	0.57
1:D:283:MET:O	1:D:287:VAL:HG23	2.03	0.57
1:E:146:ILE:HG22	1:F:462:GLN:HG2	1.86	0.57
1:H:132:TYR:OH	1:H:477:LEU:HA	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:462:GLN:HG2	1:J:146:ILE:HG22	1.86	0.57
1:E:389:VAL:HB	1:E:408:LEU:HG	1.87	0.57
1:J:132:TYR:OH	1:J:477:LEU:HA	2.04	0.57
1:D:405:MET:CE	1:D:407:ILE:HD11	2.34	0.57
1:H:377:ARG:HH11	1:H:377:ARG:HG3	1.69	0.57
1:G:302:CYS:SG	1:G:427:LEU:HD21	2.45	0.57
1:A:356:TYR:CD2	1:A:400:ILE:HG12	2.40	0.57
1:G:462:GLN:HE21	1:G:462:GLN:H	1.51	0.56
1:K:196:GLN:H	1:K:196:GLN:NE2	2.01	0.56
1:I:36:THR:CB	1:I:50:GLN:HG3	2.35	0.56
1:I:356:TYR:CD2	1:I:400:ILE:HG12	2.40	0.56
1:I:196:GLN:NE2	1:I:196:GLN:H	2.02	0.56
1:I:60:ASP:O	1:I:64:LYS:HG2	2.06	0.56
1:F:359:THR:O	1:F:363:GLU:HG3	2.05	0.56
1:D:395:ILE:HG12	5:D:3886:HOH:O	2.06	0.56
1:J:167:PRO:HD3	1:J:244:THR:HB	1.87	0.56
1:I:36:THR:HB	1:I:50:GLN:HG3	1.88	0.56
1:J:449:GLY:HA3	1:J:466:GLY:O	2.05	0.56
1:F:464:PRO:HG3	1:F:480:TYR:CD1	2.41	0.56
1:B:358:ASN:O	1:B:362:GLN:HG2	2.06	0.56
1:C:462:GLN:NE2	5:C:3931:HOH:O	2.39	0.56
1:D:361:LYS:HE2	4:D:3126:EDO:H21	1.88	0.56
1:D:196:GLN:H	1:D:196:GLN:NE2	2.03	0.55
1:J:272:LYS:HA	1:J:306:SER:OG	2.06	0.55
1:I:356:TYR:CG	1:I:400:ILE:HG12	2.41	0.55
1:I:385:VAL:HG13	1:I:405:MET:HE2	1.86	0.55
1:A:311:GLN:HG2	5:A:3752:HOH:O	2.07	0.55
1:I:135:TRP:CE2	1:K:138:LYS:HD3	2.42	0.55
1:J:12:ASN:O	1:J:15:PRO:HD3	2.06	0.55
1:C:489:LYS:HB2	1:D:468:TYR:OH	2.07	0.55
1:E:466:GLY:HA3	1:E:475:ARG:HD3	1.88	0.55
1:I:36:THR:OG1	1:I:50:GLN:HG3	2.07	0.55
1:H:464:PRO:HG3	1:H:480:TYR:CD1	2.41	0.55
1:F:46:GLU:HB2	4:F:3121:EDO:H21	1.89	0.54
1:K:462:GLN:NE2	1:K:462:GLN:H	2.04	0.54
1:L:235:HIS:HB3	1:L:238:VAL:HG23	1.89	0.54
1:A:172:LEU:HD21	1:A:200:THR:HB	1.88	0.54
1:E:167:PRO:HD3	1:E:244:THR:HB	1.87	0.54
1:K:23:PHE:CZ	1:K:26:ASN:HA	2.43	0.54
1:L:320:GLU:HB2	5:L:3748:HOH:O	2.08	0.54
1:A:167:PRO:HD3	1:A:244:THR:HB	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:294:LEU:HD12	1:B:306:SER:HA	1.90	0.54
1:L:389:VAL:HG13	1:L:393:MET:CE	2.38	0.54
1:E:459:PHE:HE2	1:E:465:PHE:CE1	2.26	0.54
1:L:11:PRO:HB3	1:L:114:TYR:CZ	2.42	0.54
1:E:23:PHE:CZ	1:E:26:ASN:HA	2.43	0.54
1:G:131:TYR:CE1	1:G:462:GLN:HB3	2.43	0.54
1:F:124:MET:HE3	1:F:173:LEU:CD2	2.38	0.54
1:G:449:GLY:HA3	1:G:466:GLY:O	2.06	0.54
1:B:449:GLY:HA3	1:B:466:GLY:O	2.08	0.54
1:G:405:MET:HE2	1:G:407:ILE:HD11	1.90	0.54
1:C:400:ILE:HD11	1:C:404:VAL:HB	1.90	0.54
1:A:254:GLN:HB2	1:A:267:LEU:HD11	1.90	0.54
1:I:283:MET:O	1:I:287:VAL:HG23	2.08	0.54
1:G:464:PRO:HG3	1:G:480:TYR:CD1	2.42	0.54
1:E:377:ARG:HG3	1:E:377:ARG:NH1	2.23	0.53
1:H:356:TYR:CD2	1:H:400:ILE:HG12	2.42	0.53
1:G:124:MET:HE3	1:G:173:LEU:CD2	2.36	0.53
1:E:124:MET:HE3	1:E:173:LEU:CD2	2.38	0.53
1:B:464:PRO:HG3	1:B:480:TYR:HD1	1.73	0.53
1:L:288:GLU:HG2	5:L:3744:HOH:O	2.07	0.53
1:E:196:GLN:H	1:E:196:GLN:NE2	2.02	0.53
1:C:155:ARG:HD2	4:C:3113:EDO:O2	2.09	0.53
1:H:167:PRO:HD3	1:H:244:THR:HB	1.90	0.53
1:D:78:ARG:HH11	1:D:78:ARG:HG2	1.73	0.53
1:G:294:LEU:HD22	1:G:405:MET:HB2	1.90	0.53
1:A:57:GLU:HG3	5:A:3762:HOH:O	2.09	0.53
1:G:172:LEU:HD21	1:G:200:THR:HB	1.91	0.53
1:F:264:ARG:HH12	1:F:484:ALA:HA	1.73	0.53
1:E:22:ILE:HD13	1:E:222:PRO:HD2	1.90	0.53
1:H:322:SER:HB3	1:H:405:MET:HE3	1.91	0.53
1:G:347:GLU:HG2	1:G:351:LYS:HE2	1.90	0.53
1:F:11:PRO:HB3	1:F:114:TYR:CZ	2.44	0.53
1:G:459:PHE:HE2	1:G:465:PHE:CE1	2.27	0.53
1:B:196:GLN:H	1:B:196:GLN:NE2	2.05	0.53
1:L:356:TYR:CD2	1:L:400:ILE:HG12	2.43	0.52
1:A:279:SER:N	1:A:311:GLN:HE21	2.06	0.52
1:G:462:GLN:NE2	5:G:3944:HOH:O	2.42	0.52
1:H:294:LEU:HD13	1:H:405:MET:HA	1.91	0.52
1:A:274:PRO:HG3	1:A:307:ARG:NH2	2.24	0.52
1:E:86:ARG:HG3	5:E:3791:HOH:O	2.08	0.52
1:L:132:TYR:OH	1:L:477:LEU:HA	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:22:ILE:HD13	1:J:222:PRO:HD2	1.91	0.52
1:L:283:MET:O	1:L:287:VAL:HG23	2.09	0.52
1:I:390:GLN:H	1:I:393:MET:HE3	1.75	0.52
1:K:443:SER:HA	1:K:451:VAL:HG11	1.92	0.52
1:C:205:ALA:HB2	1:C:220:ILE:HD12	1.92	0.51
1:G:358:ASN:HB2	5:G:3991:HOH:O	2.11	0.51
1:L:459:PHE:HE2	1:L:465:PHE:CE1	2.28	0.51
1:E:132:TYR:OH	1:E:477:LEU:HA	2.11	0.51
1:I:395:ILE:HG23	5:I:3843:HOH:O	2.11	0.51
1:C:356:TYR:CE1	1:C:400:ILE:HG22	2.45	0.51
1:F:459:PHE:HE2	1:F:465:PHE:CE1	2.29	0.51
1:I:413:ILE:O	1:I:417:VAL:HG23	2.10	0.51
1:K:459:PHE:HE2	1:K:465:PHE:CD1	2.29	0.51
1:A:405:MET:CE	1:A:407:ILE:HD11	2.40	0.51
1:A:224:PHE:HB2	1:A:227:THR:OG1	2.11	0.51
1:D:462:GLN:HE21	1:D:462:GLN:H	1.59	0.51
1:H:385:VAL:HG13	1:H:405:MET:HE2	1.92	0.51
1:K:459:PHE:HE2	1:K:465:PHE:CE1	2.28	0.51
1:C:464:PRO:HG3	1:C:480:TYR:CD1	2.46	0.51
1:H:131:TYR:CE1	1:H:462:GLN:HB3	2.46	0.50
1:J:322:SER:HB3	1:J:405:MET:HE3	1.92	0.50
1:C:18:PHE:HE1	4:C:3115:EDO:H22	1.76	0.50
1:F:264:ARG:NH1	1:F:484:ALA:HA	2.26	0.50
1:L:443:SER:HA	1:L:451:VAL:HG11	1.93	0.50
1:G:11:PRO:HB3	1:G:114:TYR:CZ	2.46	0.50
1:I:192:LYS:HB2	1:I:232:ILE:HD12	1.94	0.50
1:I:443:SER:HA	1:I:451:VAL:HG11	1.92	0.50
1:L:413:ILE:O	1:L:417:VAL:HG23	2.11	0.50
1:L:307:ARG:NH2	1:L:420:ALA:HA	2.27	0.50
1:G:311:GLN:HG2	5:G:3715:HOH:O	2.11	0.50
1:E:449:GLY:HA3	1:E:466:GLY:O	2.12	0.50
1:H:36:THR:HB	1:H:50:GLN:HG3	1.93	0.50
1:G:405:MET:CE	1:G:407:ILE:HD11	2.41	0.50
1:K:462:GLN:NE2	5:K:3787:HOH:O	2.45	0.50
1:G:23:PHE:CZ	1:G:26:ASN:HA	2.46	0.50
1:H:172:LEU:HD21	1:H:200:THR:HB	1.93	0.50
1:E:21:GLN:HB3	1:E:29:HIS:O	2.11	0.50
1:K:12:ASN:O	1:K:15:PRO:HD3	2.12	0.50
1:G:302:CYS:SG	5:G:3970:HOH:O	2.59	0.50
1:C:356:TYR:CD1	1:C:400:ILE:HG22	2.47	0.50
1:L:113:PRO:HB2	1:L:116:ILE:HG12	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:496:PRO:HB2	5:L:3787:HOH:O	2.12	0.50
1:A:251:ARG:O	1:A:255:VAL:HG23	2.12	0.49
1:I:32:VAL:HG23	1:I:58:ASP:OD1	2.12	0.49
1:G:356:TYR:CG	1:G:400:ILE:HG22	2.47	0.49
1:D:361:LYS:NZ	4:D:3126:EDO:H22	2.27	0.49
1:G:34:ARG:HG2	5:G:3948:HOH:O	2.12	0.49
1:B:298:GLN:HA	5:B:3954:HOH:O	2.12	0.49
1:D:22:ILE:HD13	1:D:222:PRO:HD2	1.94	0.49
1:C:11:PRO:HB3	1:C:114:TYR:CZ	2.47	0.49
1:H:11:PRO:HB3	1:H:114:TYR:CZ	2.47	0.49
1:D:356:TYR:CD2	1:D:400:ILE:HG12	2.47	0.49
1:H:462:GLN:H	1:H:462:GLN:CD	2.16	0.49
1:A:38:PRO:HB3	1:A:50:GLN:HE22	1.76	0.49
1:I:464:PRO:HD3	1:J:144:ILE:CD1	2.43	0.49
1:C:131:TYR:CE1	1:C:462:GLN:HB3	2.48	0.49
1:K:385:VAL:HG13	1:K:405:MET:HE2	1.93	0.49
1:C:18:PHE:CE1	4:C:3115:EDO:H22	2.48	0.49
1:E:476:GLU:HA	5:E:3839:HOH:O	2.13	0.49
1:B:356:TYR:CD2	1:B:400:ILE:HG12	2.48	0.49
1:K:196:GLN:HE21	1:K:196:GLN:N	2.05	0.49
1:J:124:MET:HE3	1:J:173:LEU:CD2	2.41	0.49
1:I:146:ILE:HG12	3:I:3107:GAI:HN31	1.78	0.49
1:D:279:SER:HA	1:D:314:ILE:HD13	1.93	0.49
1:F:131:TYR:CE1	1:F:462:GLN:HB3	2.48	0.49
1:A:294:LEU:HD13	1:A:405:MET:HA	1.95	0.49
1:A:146:ILE:HG12	3:A:3101:GAI:N3	2.28	0.49
1:C:15:PRO:HD2	1:C:108:LEU:HD22	1.95	0.49
1:D:356:TYR:CG	1:D:400:ILE:HG12	2.48	0.49
1:F:132:TYR:OH	1:F:477:LEU:HA	2.12	0.49
1:I:268:GLU:HA	5:I:3747:HOH:O	2.13	0.49
1:B:206:ASN:ND2	5:B:3986:HOH:O	2.45	0.49
1:I:124:MET:HE3	1:I:173:LEU:CD2	2.41	0.48
1:E:41:ASN:HD21	1:E:43:SER:HB2	1.78	0.48
1:H:36:THR:CB	1:H:50:GLN:HG3	2.43	0.48
1:D:391:ASP:OD2	1:D:419:ARG:HD2	2.13	0.48
1:I:196:GLN:HE21	1:I:196:GLN:N	2.06	0.48
1:J:385:VAL:HG13	1:J:405:MET:HE2	1.95	0.48
1:B:356:TYR:CE2	1:B:400:ILE:HG12	2.48	0.48
1:G:11:PRO:HB3	1:G:114:TYR:CE1	2.48	0.48
1:L:156:HIS:HB3	1:L:486:THR:HG21	1.95	0.48
1:C:196:GLN:HE21	1:C:196:GLN:N	1.99	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:120:VAL:HG12	1:F:124:MET:HE1	1.95	0.48
1:F:36:THR:CB	1:F:50:GLN:HG3	2.43	0.48
1:A:107:THR:HG23	1:A:334:PRO:HB2	1.96	0.48
1:A:315:TYR:O	1:A:319:VAL:HG23	2.14	0.48
1:E:356:TYR:CG	1:E:400:ILE:HG12	2.48	0.48
1:D:361:LYS:CE	4:D:3126:EDO:H21	2.44	0.48
1:G:464:PRO:HG3	1:G:480:TYR:HD1	1.79	0.48
1:F:32:VAL:HG11	1:F:57:GLU:OE2	2.14	0.48
1:C:302:CYS:SG	1:C:427:LEU:HD21	2.54	0.47
1:C:11:PRO:HB3	1:C:114:TYR:CE1	2.49	0.47
1:I:274:PRO:HG3	1:I:307:ARG:NH2	2.29	0.47
1:C:390:GLN:NE2	1:C:392:GLY:H	2.12	0.47
1:D:124:MET:HE3	1:D:173:LEU:CD2	2.42	0.47
1:F:120:VAL:HG12	1:F:124:MET:CE	2.45	0.47
1:D:302:CYS:SG	1:D:427:LEU:HD21	2.55	0.47
1:E:464:PRO:HG3	1:E:480:TYR:HD1	1.80	0.47
1:I:146:ILE:HG13	1:J:460:GLY:HA3	1.97	0.47
1:H:408:LEU:N	1:H:408:LEU:HD12	2.29	0.47
1:D:464:PRO:HG3	1:D:480:TYR:HD1	1.79	0.47
1:C:21:GLN:HB3	1:C:29:HIS:O	2.15	0.47
1:K:353:ILE:HG23	1:K:400:ILE:HD12	1.96	0.47
1:F:462:GLN:H	1:F:462:GLN:CD	2.18	0.47
1:L:294:LEU:CD1	1:L:306:SER:HA	2.41	0.47
1:C:270:GLY:HA3	1:C:302:CYS:HB2	1.95	0.47
1:E:462:GLN:H	1:E:462:GLN:NE2	2.12	0.47
1:B:294:LEU:HD22	1:B:405:MET:HB2	1.97	0.47
1:H:356:TYR:CE2	1:H:400:ILE:HG12	2.48	0.47
1:E:272:LYS:HA	1:E:306:SER:OG	2.14	0.47
1:L:192:LYS:HB2	1:L:232:ILE:HD12	1.97	0.47
1:H:205:ALA:HB2	1:H:220:ILE:HD12	1.97	0.47
1:I:431:VAL:HG21	1:I:442:LEU:HB3	1.96	0.47
1:K:410:PHE:CD1	1:K:416:VAL:HB	2.50	0.47
1:E:155:ARG:HD2	5:F:4007:HOH:O	2.15	0.47
1:I:241:VAL:CG1	1:I:265:VAL:HG22	2.45	0.47
1:F:174:MET:CE	1:F:244:THR:HG21	2.45	0.47
1:J:135:TRP:CE2	1:L:138:LYS:HD3	2.50	0.47
1:J:444:GLN:HB3	5:J:3751:HOH:O	2.15	0.47
1:E:131:TYR:CE1	1:E:462:GLN:HB3	2.51	0.46
1:J:356:TYR:CD2	1:J:400:ILE:HG12	2.50	0.46
1:A:356:TYR:CG	1:A:400:ILE:HG12	2.50	0.46
1:J:431:VAL:HG21	1:J:442:LEU:HB3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:463:SER:HA	1:B:464:PRO:HD3	1.80	0.46
1:E:36:THR:OG1	1:E:50:GLN:HG3	2.14	0.46
1:B:268:GLU:HG3	1:B:268:GLU:O	2.16	0.46
1:D:166:ILE:HD11	1:D:193:VAL:HG12	1.98	0.46
1:E:468:TYR:OH	1:F:489:LYS:HB2	2.15	0.46
1:J:405:MET:CE	1:J:407:ILE:HD11	2.46	0.46
1:I:12:ASN:O	1:I:15:PRO:HD3	2.15	0.46
1:E:124:MET:HE1	5:E:3985:HOH:O	2.15	0.46
1:J:464:PRO:HG3	1:J:480:TYR:CD1	2.50	0.46
1:C:101:TYR:CG	4:C:3115:EDO:H11	2.50	0.46
1:D:156:HIS:HB3	1:D:486:THR:HG21	1.97	0.46
1:K:464:PRO:HG3	1:K:480:TYR:CD1	2.51	0.46
1:E:11:PRO:HB3	1:E:114:TYR:CZ	2.50	0.46
1:G:32:VAL:HG21	1:G:57:GLU:HG3	1.97	0.46
1:G:132:TYR:OH	1:G:477:LEU:HA	2.15	0.46
1:K:146:ILE:HD11	3:L:3108:GAI:N3	2.30	0.46
1:L:415:GLU:HB2	5:L:3799:HOH:O	2.16	0.46
1:C:466:GLY:HA3	1:C:475:ARG:HD3	1.98	0.46
1:C:356:TYR:CD1	1:C:400:ILE:CG2	2.99	0.46
1:A:233:ALA:HB2	1:A:253:ILE:HG23	1.96	0.46
1:F:431:VAL:HG21	1:F:442:LEU:HB3	1.97	0.46
1:A:294:LEU:HD22	1:A:405:MET:HB2	1.98	0.46
1:B:462:GLN:NE2	5:B:3957:HOH:O	2.49	0.46
1:I:468:TYR:OH	1:J:489:LYS:HB2	2.16	0.46
1:B:120:VAL:HG12	1:B:124:MET:CE	2.45	0.46
1:L:120:VAL:HG12	1:L:124:MET:HE1	1.98	0.46
1:F:174:MET:HE1	1:F:244:THR:HG21	1.98	0.46
1:I:294:LEU:HD13	1:I:294:LEU:C	2.36	0.46
1:E:294:LEU:O	1:E:294:LEU:HG	2.15	0.45
1:D:464:PRO:CG	1:D:480:TYR:CD1	2.99	0.45
1:K:11:PRO:HB3	1:K:114:TYR:CZ	2.51	0.45
1:G:460:GLY:HA3	1:H:146:ILE:HG13	1.98	0.45
1:J:23:PHE:CZ	1:J:26:ASN:HA	2.51	0.45
1:B:131:TYR:CE1	1:B:462:GLN:HB3	2.51	0.45
1:A:38:PRO:HB3	1:A:50:GLN:NE2	2.30	0.45
1:E:205:ALA:HB2	1:E:220:ILE:HD12	1.99	0.45
1:B:205:ALA:HB2	1:B:220:ILE:HD12	1.98	0.45
1:H:476:GLU:O	1:H:477:LEU:HB2	2.17	0.45
1:H:449:GLY:HA3	1:H:466:GLY:O	2.16	0.45
1:J:443:SER:HA	1:J:451:VAL:HG11	1.98	0.45
1:L:283:MET:HG3	1:L:321:ARG:NH1	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:465:PHE:O	1:K:476:GLU:HB3	2.15	0.45
1:K:121:ASP:O	1:K:125:VAL:HG23	2.16	0.45
1:F:487:LYS:HZ1	3:F:3103:GAI:C	2.30	0.45
1:A:46:GLU:CB	4:A:3110:EDO:H21	2.46	0.45
1:A:23:PHE:CZ	1:A:26:ASN:HA	2.52	0.45
1:F:101:TYR:CG	4:F:3122:EDO:H11	2.52	0.45
1:D:322:SER:HB3	1:D:405:MET:HE3	1.99	0.45
1:J:356:TYR:CG	1:J:400:ILE:HG12	2.52	0.45
1:C:356:TYR:CG	1:C:400:ILE:HG21	2.51	0.45
1:H:36:THR:OG1	1:H:50:GLN:HG3	2.16	0.45
1:I:445:ALA:HB2	5:I:3756:HOH:O	2.17	0.45
1:C:272:LYS:HA	1:C:306:SER:OG	2.16	0.45
1:D:353:ILE:O	1:D:357:ILE:HG13	2.16	0.45
1:D:449:GLY:HA3	1:D:466:GLY:O	2.17	0.45
1:D:409:LYS:O	1:D:419:ARG:NH2	2.50	0.45
1:D:46:GLU:HB2	4:D:3117:EDO:H21	1.99	0.45
1:A:131:TYR:CZ	1:A:462:GLN:HA	2.52	0.45
1:A:273:SER:HA	1:A:274:PRO:HD3	1.74	0.45
1:K:167:PRO:HD3	1:K:244:THR:O	2.16	0.45
1:B:464:PRO:CG	1:B:480:TYR:CD1	2.98	0.44
1:I:463:SER:HA	1:I:464:PRO:HD3	1.85	0.44
1:H:389:VAL:HB	1:H:408:LEU:HG	1.98	0.44
1:C:214:PRO:HD2	1:C:217:VAL:HG21	1.98	0.44
1:J:46:GLU:HB2	4:J:3128:EDO:H21	1.98	0.44
1:E:302:CYS:HB3	5:E:3850:HOH:O	2.17	0.44
1:L:15:PRO:HD2	1:L:108:LEU:HD22	1.99	0.44
1:A:32:VAL:HG11	1:A:57:GLU:OE2	2.17	0.44
1:G:21:GLN:HB3	1:G:29:HIS:O	2.17	0.44
1:D:443:SER:HA	1:D:451:VAL:HG11	1.99	0.44
1:J:161:VAL:HA	1:J:188:VAL:HG23	1.99	0.44
1:F:311:GLN:HG2	5:F:3780:HOH:O	2.17	0.44
1:A:138:LYS:HD3	1:C:135:TRP:CE2	2.52	0.44
1:K:319:VAL:O	1:K:323:VAL:HG23	2.17	0.44
1:B:192:LYS:HB2	1:B:232:ILE:HD12	1.99	0.44
1:L:243:PHE:HB3	1:L:267:LEU:HD23	1.98	0.44
1:G:327:LYS:HE3	1:G:369:CYS:HB3	1.99	0.44
1:B:294:LEU:CD1	1:B:306:SER:HA	2.47	0.44
1:E:36:THR:CB	1:E:50:GLN:HG3	2.47	0.44
1:L:33:SER:O	1:L:34:ARG:HB2	2.17	0.44
1:B:459:PHE:HE2	1:B:465:PHE:CD1	2.35	0.44
1:K:174:MET:CE	1:K:177:TRP:CZ3	3.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:361:LYS:HZ3	4:D:3126:EDO:H22	1.82	0.44
1:D:243:PHE:HB3	1:D:267:LEU:HD23	1.98	0.44
1:H:23:PHE:CZ	1:H:26:ASN:HA	2.53	0.44
1:C:46:GLU:HB2	4:C:3114:EDO:H21	2.00	0.44
1:L:310:VAL:HG21	1:L:318:PHE:CD2	2.52	0.44
1:E:41:ASN:C	1:E:41:ASN:ND2	2.69	0.44
1:E:273:SER:HA	1:E:274:PRO:HD3	1.78	0.44
1:K:192:LYS:HB2	1:K:232:ILE:HD12	2.00	0.44
1:I:453:VAL:HB	1:J:493:VAL:HG13	2.00	0.44
1:E:106:GLU:O	1:E:110:ASN:HB3	2.18	0.44
1:H:272:LYS:HA	1:H:306:SER:OG	2.18	0.44
1:K:294:LEU:HD12	1:K:306:SER:HA	2.00	0.44
1:K:39:THR:HG23	1:K:48:ILE:HB	1.99	0.44
1:H:64:LYS:HE2	5:H:3932:HOH:O	2.18	0.44
1:L:459:PHE:O	3:L:3108:GAI:N2	2.49	0.44
1:G:385:VAL:HG13	1:G:405:MET:HE2	2.00	0.44
1:I:310:VAL:HG21	1:I:318:PHE:CD2	2.53	0.44
1:I:193:VAL:HG11	1:I:201:ALA:CB	2.48	0.44
1:A:349:GLN:O	1:A:353:ILE:HG13	2.18	0.44
1:B:243:PHE:HB3	1:B:267:LEU:HD23	2.00	0.44
1:B:400:ILE:H	1:B:400:ILE:HG13	1.71	0.44
1:C:172:LEU:HD21	1:C:200:THR:HB	1.99	0.44
1:H:333:ASN:HA	1:H:334:PRO:HD2	1.85	0.44
1:D:315:TYR:O	1:D:319:VAL:HG23	2.18	0.44
1:D:188:VAL:HG12	1:D:217:VAL:HA	2.00	0.44
1:L:431:VAL:HG21	1:L:442:LEU:HB3	1.99	0.44
1:F:159:VAL:HG21	1:F:264:ARG:HH21	1.83	0.43
1:I:146:ILE:HG12	3:I:3107:GAI:N3	2.33	0.43
1:D:161:VAL:HA	1:D:188:VAL:HG23	1.99	0.43
1:I:158:PRO:HG3	1:I:185:THR:O	2.18	0.43
1:L:315:TYR:CE2	1:L:409:LYS:HB2	2.53	0.43
1:L:131:TYR:CE1	1:L:462:GLN:HB3	2.52	0.43
1:A:174:MET:HE2	1:A:177:TRP:CE3	2.53	0.43
1:I:292:PHE:HE1	1:I:457:ASP:HB2	1.82	0.43
1:B:496:PRO:HG2	1:D:441:TYR:HB2	2.00	0.43
1:D:131:TYR:CE1	1:D:462:GLN:HB3	2.53	0.43
1:H:462:GLN:NE2	5:H:3866:HOH:O	2.52	0.43
1:L:356:TYR:CE2	1:L:400:ILE:HG12	2.54	0.43
1:A:319:VAL:O	1:A:323:VAL:HG23	2.19	0.43
1:L:21:GLN:HB3	1:L:29:HIS:O	2.17	0.43
1:J:235:HIS:HB3	1:J:238:VAL:HG23	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:283:MET:O	1:H:287:VAL:HG23	2.17	0.43
1:K:272:LYS:HA	1:K:306:SER:OG	2.18	0.43
1:E:356:TYR:CE2	1:E:400:ILE:HG12	2.54	0.43
1:J:464:PRO:HA	1:J:476:GLU:O	2.18	0.43
1:D:59:VAL:O	1:D:63:VAL:HG23	2.18	0.43
1:B:158:PRO:HG3	1:B:185:THR:O	2.19	0.43
1:A:443:SER:HA	1:A:451:VAL:HG11	2.00	0.43
1:H:464:PRO:CG	1:H:480:TYR:CD1	3.01	0.43
1:J:405:MET:HE3	1:J:407:ILE:HD11	1.99	0.43
1:C:449:GLY:HA3	1:C:466:GLY:O	2.18	0.43
1:I:205:ALA:HB2	1:I:220:ILE:HD12	2.00	0.43
1:F:59:VAL:O	1:F:63:VAL:HG23	2.18	0.43
1:I:131:TYR:CE1	1:I:462:GLN:HB3	2.54	0.43
1:B:124:MET:HE3	1:B:173:LEU:CD2	2.48	0.43
1:F:27:GLU:HG3	5:F:3816:HOH:O	2.18	0.43
1:F:272:LYS:HA	1:F:306:SER:OG	2.18	0.43
1:I:167:PRO:HD3	1:I:244:THR:HB	2.00	0.43
1:C:13:GLN:NE2	5:C:3894:HOH:O	2.51	0.43
1:K:146:ILE:HG12	3:L:3108:GAI:N3	2.30	0.43
1:E:36:THR:HB	1:E:50:GLN:HG3	2.00	0.43
1:J:11:PRO:HB3	1:J:114:TYR:CZ	2.54	0.43
1:B:156:HIS:CD2	1:B:488:VAL:HG22	2.54	0.43
1:D:377:ARG:HG3	1:D:377:ARG:HH11	1.83	0.43
1:A:294:LEU:CD1	1:A:306:SER:HA	2.46	0.43
1:F:464:PRO:CG	1:F:480:TYR:CD1	3.01	0.43
1:I:154:THR:HG21	1:J:464:PRO:HG2	2.00	0.43
1:D:192:LYS:HD3	1:D:193:VAL:O	2.19	0.43
1:E:302:CYS:SG	1:E:427:LEU:HD21	2.59	0.43
1:H:315:TYR:CE1	1:H:319:VAL:HG21	2.54	0.43
1:K:404:VAL:HG12	1:K:406:GLN:OE1	2.19	0.43
1:C:294:LEU:HG	1:C:294:LEU:O	2.10	0.43
1:I:464:PRO:HG3	1:I:480:TYR:CD1	2.54	0.43
1:E:8:VAL:HA	1:E:9:PRO:HD3	1.80	0.43
1:C:431:VAL:HG21	1:C:442:LEU:HB3	2.00	0.43
1:K:146:ILE:HG13	1:L:460:GLY:HA3	2.01	0.42
1:D:294:LEU:CD1	1:D:405:MET:HA	2.47	0.42
1:K:405:MET:HE3	1:K:407:ILE:HD11	1.99	0.42
1:L:349:GLN:O	1:L:353:ILE:HG13	2.19	0.42
1:F:192:LYS:NZ	5:F:3778:HOH:O	2.52	0.42
1:A:235:HIS:HB3	1:A:238:VAL:HG23	2.00	0.42
1:E:356:TYR:HB2	1:E:400:ILE:HG21	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:373:ILE:HG22	1:K:375:ALA:H	1.84	0.42
1:D:34:ARG:HH11	1:D:34:ARG:HB3	1.84	0.42
1:D:462:GLN:CD	1:D:462:GLN:H	2.22	0.42
1:B:462:GLN:HE21	1:B:462:GLN:H	1.64	0.42
1:K:132:TYR:OH	1:K:477:LEU:HA	2.19	0.42
1:H:377:ARG:HG3	1:H:377:ARG:NH1	2.34	0.42
1:B:466:GLY:HA3	1:B:475:ARG:HD3	2.02	0.42
1:D:410:PHE:CD1	1:D:416:VAL:HB	2.54	0.42
1:I:349:GLN:O	1:I:353:ILE:HG13	2.18	0.42
1:F:241:VAL:CG1	1:F:265:VAL:HG22	2.49	0.42
1:D:373:ILE:HG22	1:D:375:ALA:H	1.85	0.42
1:J:347:GLU:HG2	1:J:351:LYS:HE2	2.00	0.42
1:D:158:PRO:HG3	1:D:185:THR:O	2.19	0.42
1:K:345:VAL:HG13	1:K:346:ASP:N	2.35	0.42
1:H:322:SER:HB3	1:H:405:MET:CE	2.49	0.42
1:H:107:THR:HG23	1:H:334:PRO:HB2	2.02	0.42
1:L:264:ARG:NH1	1:L:484:ALA:O	2.52	0.42
1:E:468:TYR:HE1	1:E:475:ARG:HH21	1.67	0.42
1:C:495:VAL:C	1:C:496:PRO:O	2.57	0.42
1:J:59:VAL:O	1:J:63:VAL:HG23	2.20	0.42
1:A:8:VAL:HA	1:A:9:PRO:HD3	1.93	0.42
1:H:372:GLY:O	1:H:382:GLN:HG3	2.19	0.42
1:K:322:SER:HB3	1:K:405:MET:HE3	2.00	0.42
1:J:464:PRO:HG3	1:J:480:TYR:HD1	1.85	0.42
1:C:400:ILE:HG21	1:C:400:ILE:HD13	1.78	0.42
1:D:21:GLN:HB3	1:D:29:HIS:O	2.19	0.42
1:I:214:PRO:HA	1:I:215:PRO:HD3	1.94	0.42
1:I:315:TYR:CE1	1:I:319:VAL:HG21	2.55	0.42
1:J:158:PRO:HG3	1:J:185:THR:O	2.20	0.42
1:L:294:LEU:CD1	1:L:405:MET:HA	2.50	0.42
1:H:405:MET:HE3	1:H:407:ILE:HD11	2.01	0.42
1:I:461:ALA:N	1:I:462:GLN:NE2	2.68	0.42
1:D:361:LYS:CE	4:D:3126:EDO:C2	2.98	0.42
1:I:405:MET:HE3	1:I:407:ILE:HD11	2.02	0.42
1:F:476:GLU:O	1:F:477:LEU:HB2	2.19	0.42
1:G:46:GLU:HB2	4:G:3124:EDO:H21	2.01	0.42
1:F:463:SER:HA	1:F:464:PRO:HD3	1.79	0.41
1:A:353:ILE:O	1:A:357:ILE:HG13	2.20	0.41
1:F:12:ASN:O	1:F:15:PRO:HD3	2.19	0.41
1:B:132:TYR:OH	1:B:477:LEU:HA	2.19	0.41
1:L:463:SER:HA	1:L:464:PRO:HD3	1.87	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:164:GLN:CD	1:A:178:LYS:HB3	2.39	0.41
1:D:121:ASP:O	1:D:125:VAL:HG23	2.21	0.41
1:L:60:ASP:O	1:L:64:LYS:HG2	2.19	0.41
1:E:464:PRO:CG	1:E:480:TYR:CD1	3.02	0.41
1:J:356:TYR:HE2	1:J:404:VAL:HG11	1.84	0.41
1:J:107:THR:HG23	1:J:334:PRO:HB2	2.01	0.41
1:B:422:ASN:HB3	5:B:3794:HOH:O	2.20	0.41
1:K:431:VAL:HG21	1:K:442:LEU:HB3	2.02	0.41
1:I:273:SER:HA	1:I:274:PRO:HD3	1.80	0.41
1:D:235:HIS:HB3	1:D:238:VAL:HG23	2.02	0.41
1:B:389:VAL:HB	1:B:408:LEU:HG	2.02	0.41
1:I:57:GLU:H	1:I:57:GLU:CD	2.24	0.41
1:A:322:SER:HB3	1:A:405:MET:CE	2.51	0.41
1:J:462:GLN:H	1:J:462:GLN:CD	2.22	0.41
1:C:154:THR:HA	1:C:489:LYS:O	2.20	0.41
1:B:459:PHE:HE2	1:B:465:PHE:CE1	2.39	0.41
1:B:315:TYR:O	1:B:319:VAL:HG23	2.20	0.41
1:F:8:VAL:HA	1:F:9:PRO:HD3	1.94	0.41
1:K:302:CYS:HA	1:K:401:PHE:HE2	1.85	0.41
1:E:112:LYS:HB3	1:E:112:LYS:HE2	1.92	0.41
1:K:347:GLU:HG2	1:K:351:LYS:HE2	2.03	0.41
1:F:196:GLN:N	1:F:196:GLN:HE21	2.09	0.41
1:I:144:ILE:CD1	1:J:464:PRO:HD3	2.51	0.41
1:D:192:LYS:HZ3	1:D:225:GLY:HA2	1.85	0.41
1:H:12:ASN:O	1:H:15:PRO:HD3	2.21	0.41
1:L:462:GLN:H	1:L:462:GLN:HE21	1.64	0.41
1:E:463:SER:HA	1:E:464:PRO:HD3	1.85	0.41
1:A:167:PRO:HG2	1:A:174:MET:HG3	2.02	0.41
1:J:356:TYR:CE1	1:J:395:ILE:HG22	2.55	0.41
1:D:192:LYS:NZ	1:D:225:GLY:HA2	2.35	0.41
1:H:273:SER:HA	1:H:274:PRO:HD3	1.83	0.41
1:A:408:LEU:HD12	1:A:408:LEU:N	2.35	0.41
1:A:131:TYR:HE1	1:A:462:GLN:HG3	1.79	0.41
1:L:146:ILE:HD11	1:L:150:PHE:HB2	2.02	0.41
1:A:347:GLU:O	1:A:350:PHE:HB3	2.20	0.41
1:D:418:GLY:HA3	5:D:3775:HOH:O	2.20	0.41
1:L:345:VAL:HG13	1:L:346:ASP:N	2.35	0.41
1:F:46:GLU:CB	4:F:3121:EDO:H21	2.50	0.41
1:K:408:LEU:HD12	1:K:408:LEU:N	2.36	0.41
1:I:466:GLY:HA3	1:I:475:ARG:CD	2.38	0.41
1:F:294:LEU:HG	1:F:294:LEU:O	2.18	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:356:TYR:CG	1:H:400:ILE:HG12	2.55	0.41
1:B:356:TYR:CG	1:B:400:ILE:HG12	2.56	0.41
1:F:177:TRP:CD1	1:F:477:LEU:HD21	2.56	0.41
1:E:155:ARG:NH2	1:F:444:GLN:HG3	2.36	0.41
1:C:23:PHE:CZ	1:C:26:ASN:HA	2.56	0.41
1:K:235:HIS:HB3	1:K:238:VAL:HG23	2.01	0.41
1:D:23:PHE:CZ	1:D:26:ASN:HA	2.56	0.41
1:A:11:PRO:HB3	1:A:114:TYR:CZ	2.56	0.41
1:C:490:THR:O	1:D:450:THR:HA	2.21	0.41
1:L:166:ILE:HD11	5:L:3726:HOH:O	2.21	0.41
1:L:167:PRO:HD3	1:L:244:THR:HB	2.02	0.41
1:L:294:LEU:HD13	1:L:405:MET:HA	2.03	0.41
4:A:3116:EDO:C1	1:C:444:GLN:HG3	2.51	0.41
1:K:372:GLY:O	1:K:382:GLN:HG3	2.21	0.41
1:D:34:ARG:NH1	1:D:34:ARG:HB3	2.37	0.40
1:L:285:TRP:CH2	1:L:289:GLN:NE2	2.90	0.40
1:L:377:ARG:HG3	1:L:377:ARG:HH11	1.86	0.40
1:C:132:TYR:OH	1:C:477:LEU:HA	2.21	0.40
1:F:23:PHE:CZ	1:F:26:ASN:HA	2.57	0.40
1:C:102:LEU:HD23	1:C:102:LEU:HA	1.91	0.40
1:L:333:ASN:HA	1:L:334:PRO:HD2	1.94	0.40
1:A:146:ILE:HG12	3:A:3101:GAI:HN31	1.87	0.40
1:H:410:PHE:CD1	1:H:416:VAL:HB	2.57	0.40
1:G:121:ASP:O	1:G:125:VAL:HG23	2.20	0.40
1:L:193:VAL:HG11	1:L:201:ALA:CB	2.51	0.40
1:C:166:ILE:HD11	5:C:3707:HOH:O	2.20	0.40
1:D:167:PRO:HG2	1:D:174:MET:HG3	2.01	0.40
1:I:112:LYS:HE2	1:I:112:LYS:HB3	1.96	0.40
1:H:406:GLN:N	1:H:406:GLN:NE2	2.69	0.40
1:K:279:SER:HA	1:K:314:ILE:HD13	2.04	0.40
1:I:462:GLN:HE21	1:I:462:GLN:H	1.67	0.40
1:F:459:PHE:HE2	1:F:465:PHE:CD1	2.40	0.40
1:C:463:SER:HA	1:C:464:PRO:HD3	1.79	0.40
1:G:34:ARG:HD2	1:K:16:GLU:OE2	2.21	0.40
1:L:192:LYS:HB2	1:L:232:ILE:CD1	2.51	0.40
1:D:464:PRO:HG3	1:D:480:TYR:CD1	2.56	0.40
1:E:274:PRO:HG3	1:E:307:ARG:NH2	2.36	0.40
1:I:72:LEU:HD12	1:I:72:LEU:HA	1.86	0.40

There are no symmetry-related clashes.



## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	461/500 (92%)	451 (98%)	10 (2%)	0	100	100
1	B	457/500 (91%)	449 (98%)	7 (2%)	1 (0%)	52	53
1	C	460/500 (92%)	451 (98%)	8 (2%)	1 (0%)	52	53
1	D	460/500 (92%)	448 (97%)	12 (3%)	0	100	100
1	E	461/500 (92%)	449 (97%)	12 (3%)	0	100	100
1	F	460/500 (92%)	451 (98%)	7 (2%)	2 (0%)	39	37
1	G	461/500 (92%)	452 (98%)	8 (2%)	1 (0%)	52	53
1	H	461/500 (92%)	452 (98%)	8 (2%)	1 (0%)	52	53
1	I	460/500 (92%)	448 (97%)	10 (2%)	2 (0%)	39	37
1	J	459/500 (92%)	448 (98%)	10 (2%)	1 (0%)	52	53
1	K	448/500 (90%)	437 (98%)	10 (2%)	1 (0%)	52	53
1	L	454/500 (91%)	443 (98%)	10 (2%)	1 (0%)	52	53
All	All	5502/6000 (92%)	5379 (98%)	112 (2%)	11 (0%)	52	53

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	496	PRO
1	F	466	GLY
1	I	466	GLY
1	J	496	PRO
1	C	496	PRO
1	K	496	PRO
1	L	496	PRO
1	F	496	PRO
1	I	496	PRO
1	G	496	PRO
1	H	496	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	381/402 (95%)	375 (98%)	6 (2%)	70	76
1	B	376/402 (94%)	368 (98%)	8 (2%)	61	66
1	C	379/402 (94%)	370 (98%)	9 (2%)	57	61
1	D	379/402 (94%)	373 (98%)	6 (2%)	70	76
1	E	380/402 (94%)	373 (98%)	7 (2%)	66	72
1	F	379/402 (94%)	372 (98%)	7 (2%)	66	72
1	G	380/402 (94%)	376 (99%)	4 (1%)	80	85
1	H	380/402 (94%)	374 (98%)	6 (2%)	70	76
1	I	379/402 (94%)	374 (99%)	5 (1%)	76	82
1	J	378/402 (94%)	371 (98%)	7 (2%)	65	70
1	K	371/402 (92%)	365 (98%)	6 (2%)	70	76
1	L	376/402 (94%)	371 (99%)	5 (1%)	76	82
All	All	4538/4824 (94%)	4462 (98%)	76 (2%)	68	74

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	192	LYS
1	A	254	GLN
1	A	275	ASN
1	A	376	ASP
1	A	462	GLN
1	A	497	GLN
1	B	56	LYS
1	B	60	ASP
1	B	192	LYS
1	B	196	GLN
1	B	268	GLU
1	B	294	LEU
1	B	462	GLN
1	B	496	PRO

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Mol	Chain	Res	Type
1	C	122	LEU
1	C	192	LYS
1	C	196	GLN
1	C	236	GLU
1	C	275	ASN
1	C	376	ASP
1	C	377	ARG
1	C	462	GLN
1	C	496	PRO
1	D	192	LYS
1	D	196	GLN
1	D	275	ASN
1	D	358	ASN
1	D	362	GLN
1	D	462	GLN
1	E	41	ASN
1	E	122	LEU
1	E	196	GLN
1	E	264	ARG
1	E	275	ASN
1	E	294	LEU
1	E	376	ASP
1	F	34	ARG
1	F	122	LEU
1	F	192	LYS
1	F	196	GLN
1	F	400	ILE
1	F	462	GLN
1	F	496	PRO
1	G	192	LYS
1	G	294	LEU
1	G	383	PRO
1	G	462	GLN
1	H	192	LYS
1	H	196	GLN
1	H	275	ASN
1	H	376	ASP
1	H	462	GLN
1	H	463	SER
1	I	192	LYS
1	I	196	GLN
1	I	275	ASN

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Mol	Chain	Res	Type
1	I	462	GLN
1	I	463	SER
1	J	36	THR
1	J	50	GLN
1	J	192	LYS
1	J	196	GLN
1	J	294	LEU
1	J	462	GLN
1	J	496	PRO
1	K	192	LYS
1	K	196	GLN
1	K	294	LEU
1	K	325	ARG
1	K	462	GLN
1	K	496	PRO
1	L	117	SER
1	L	192	LYS
1	L	275	ASN
1	L	409	LYS
1	L	462	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (92) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	13	GLN
1	A	26	ASN
1	A	50	GLN
1	A	164	GLN
1	A	175	GLN
1	A	254	GLN
1	A	275	ASN
1	A	311	GLN
1	A	358	ASN
1	A	444	GLN
1	A	462	GLN
1	A	483	GLN
1	B	14	GLN
1	B	26	ASN
1	B	71	GLN
1	B	83	HIS
1	B	156	HIS
1	B	175	GLN

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Mol	Chain	Res	Type
1	B	196	GLN
1	B	275	ASN
1	B	358	ASN
1	B	462	GLN
1	C	13	GLN
1	C	26	ASN
1	C	50	GLN
1	C	83	HIS
1	C	175	GLN
1	C	196	GLN
1	C	275	ASN
1	C	390	GLN
1	C	422	ASN
1	C	462	GLN
1	D	13	GLN
1	D	26	ASN
1	D	83	HIS
1	D	175	GLN
1	D	196	GLN
1	D	275	ASN
1	D	462	GLN
1	E	13	GLN
1	E	41	ASN
1	E	175	GLN
1	E	196	GLN
1	E	275	ASN
1	E	462	GLN
1	F	26	ASN
1	F	50	GLN
1	F	175	GLN
1	F	196	GLN
1	F	275	ASN
1	F	462	GLN
1	G	13	GLN
1	G	26	ASN
1	G	175	GLN
1	G	275	ASN
1	G	462	GLN
1	H	26	ASN
1	H	164	GLN
1	H	175	GLN
1	H	196	GLN

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Mol	Chain	Res	Type
1	H	275	ASN
1	H	462	GLN
1	I	13	GLN
1	I	26	ASN
1	I	50	GLN
1	I	83	HIS
1	I	175	GLN
1	I	196	GLN
1	I	275	ASN
1	I	462	GLN
1	J	13	GLN
1	J	21	GLN
1	J	26	ASN
1	J	83	HIS
1	J	196	GLN
1	J	275	ASN
1	J	462	GLN
1	J	483	GLN
1	K	13	GLN
1	K	26	ASN
1	K	83	HIS
1	K	196	GLN
1	K	275	ASN
1	K	462	GLN
1	L	13	GLN
1	L	14	GLN
1	L	26	ASN
1	L	50	GLN
1	L	83	HIS
1	L	175	GLN
1	L	275	ASN
1	L	462	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

Of 41 ligands modelled in this entry, 12 are monoatomic - leaving 29 for Mogul analysis.

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	GAI	A	3101	-	0,3,3	0.00	-	0,3,3	0.00	-
4	EDO	A	3109	-	3,3,3	0.44	0	2,2,2	0.37	0
4	EDO	A	3110	-	3,3,3	0.44	0	2,2,2	0.38	0
4	EDO	A	3111	-	3,3,3	0.60	0	2,2,2	0.39	0
4	EDO	A	3116	-	3,3,3	0.63	0	2,2,2	0.15	0
4	EDO	B	3112	-	3,3,3	0.44	0	2,2,2	0.39	0
4	EDO	C	3113	-	3,3,3	0.69	0	2,2,2	0.16	0
4	EDO	C	3114	-	3,3,3	0.29	0	2,2,2	0.50	0
4	EDO	C	3115	-	3,3,3	0.41	0	2,2,2	0.55	0
4	EDO	D	3117	-	3,3,3	0.50	0	2,2,2	0.31	0
4	EDO	D	3126	-	3,3,3	0.43	0	2,2,2	0.49	0
3	GAI	E	3102	-	0,3,3	0.00	-	0,3,3	0.00	-
4	EDO	E	3118	-	3,3,3	0.46	0	2,2,2	0.36	0
3	GAI	F	3103	-	0,3,3	0.00	-	0,3,3	0.00	-
3	GAI	F	3104	-	0,3,3	0.00	-	0,3,3	0.00	-
4	EDO	F	3120	-	3,3,3	0.40	0	2,2,2	0.41	0
4	EDO	F	3121	-	3,3,3	0.44	0	2,2,2	0.47	0
4	EDO	F	3122	-	3,3,3	0.36	0	2,2,2	0.60	0
3	GAI	G	3105	-	0,3,3	0.00	-	0,3,3	0.00	-
4	EDO	G	3123	-	3,3,3	0.43	0	2,2,2	0.47	0
4	EDO	G	3124	-	3,3,3	0.49	0	2,2,2	0.32	0
4	EDO	G	3125	-	3,3,3	0.50	0	2,2,2	0.48	0
3	GAI	H	3106	-	0,3,3	0.00	-	0,3,3	0.00	-
3	GAI	I	3107	-	0,3,3	0.00	-	0,3,3	0.00	-
4	EDO	I	3119	-	3,3,3	0.34	0	2,2,2	0.65	0
4	EDO	I	3127	-	3,3,3	0.40	0	2,2,2	0.60	0
4	EDO	J	3128	-	3,3,3	0.49	0	2,2,2	0.29	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	GAI	L	3108	-	0,3,3	0.00	-	0,3,3	0.00	-
4	EDO	L	3129	-	3,3,3	0.52	0	2,2,2	0.35	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	GAI	A	3101	-	-	0/0/0/0	0/0/0/0
4	EDO	A	3109	-	-	0/1/1/1	0/0/0/0
4	EDO	A	3110	-	-	0/1/1/1	0/0/0/0
4	EDO	A	3111	-	-	0/1/1/1	0/0/0/0
4	EDO	A	3116	-	-	0/1/1/1	0/0/0/0
4	EDO	B	3112	-	-	0/1/1/1	0/0/0/0
4	EDO	C	3113	-	-	0/1/1/1	0/0/0/0
4	EDO	C	3114	-	-	0/1/1/1	0/0/0/0
4	EDO	C	3115	-	-	0/1/1/1	0/0/0/0
4	EDO	D	3117	-	-	0/1/1/1	0/0/0/0
4	EDO	D	3126	-	-	0/1/1/1	0/0/0/0
3	GAI	E	3102	-	-	0/0/0/0	0/0/0/0
4	EDO	E	3118	-	-	0/1/1/1	0/0/0/0
3	GAI	F	3103	-	-	0/0/0/0	0/0/0/0
3	GAI	F	3104	-	-	0/0/0/0	0/0/0/0
4	EDO	F	3120	-	-	0/1/1/1	0/0/0/0
4	EDO	F	3121	-	-	0/1/1/1	0/0/0/0
4	EDO	F	3122	-	-	0/1/1/1	0/0/0/0
3	GAI	G	3105	-	-	0/0/0/0	0/0/0/0
4	EDO	G	3123	-	-	0/1/1/1	0/0/0/0
4	EDO	G	3124	-	-	0/1/1/1	0/0/0/0
4	EDO	G	3125	-	-	0/1/1/1	0/0/0/0
3	GAI	H	3106	-	-	0/0/0/0	0/0/0/0
3	GAI	I	3107	-	-	0/0/0/0	0/0/0/0
4	EDO	I	3119	-	-	0/1/1/1	0/0/0/0
4	EDO	I	3127	-	-	0/1/1/1	0/0/0/0
4	EDO	J	3128	-	-	0/1/1/1	0/0/0/0
3	GAI	L	3108	-	-	0/0/0/0	0/0/0/0
4	EDO	L	3129	-	-	0/1/1/1	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.



There are no torsion outliers.

There are no ring outliers.

19 monomers are involved in 36 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	3101	GAI	2	0
4	A	3109	EDO	1	0
4	A	3110	EDO	2	0
4	A	3116	EDO	5	0
4	C	3113	EDO	1	0
4	C	3114	EDO	1	0
4	C	3115	EDO	3	0
4	D	3117	EDO	1	0
4	D	3126	EDO	5	0
3	F	3103	GAI	1	0
4	F	3120	EDO	1	0
4	F	3121	EDO	2	0
4	F	3122	EDO	1	0
4	G	3123	EDO	1	0
4	G	3124	EDO	1	0
3	I	3107	GAI	2	0
4	J	3128	EDO	1	0
3	L	3108	GAI	4	0
4	L	3129	EDO	1	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	469/500 (93%)	0.52	38 (8%) 15 20	13, 49, 74, 82	0
1	B	465/500 (93%)	0.00	12 (2%) 59 66	15, 28, 50, 92	0
1	C	468/500 (93%)	-0.09	12 (2%) 59 66	14, 27, 56, 83	0
1	D	468/500 (93%)	0.45	41 (8%) 12 17	15, 45, 71, 83	0
1	E	469/500 (93%)	-0.02	14 (2%) 54 62	12, 29, 53, 83	0
1	F	468/500 (93%)	-0.19	9 (1%) 70 75	13, 24, 48, 83	0
1	G	469/500 (93%)	-0.03	13 (2%) 56 64	15, 30, 56, 79	0
1	H	469/500 (93%)	0.04	14 (2%) 54 62	13, 33, 62, 78	0
1	I	468/500 (93%)	0.34	33 (7%) 19 26	25, 43, 74, 86	0
1	J	467/500 (93%)	0.31	20 (4%) 39 48	26, 47, 70, 87	0
1	K	456/500 (91%)	0.51	33 (7%) 18 25	34, 53, 73, 92	0
1	L	462/500 (92%)	1.32	115 (24%) 1 1	32, 64, 88, 95	0
All	All	5598/6000 (93%)	0.26	354 (6%) 23 31	12, 39, 74, 95	0

All (354) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	267	LEU	8.3
1	I	413	ILE	8.2
1	L	267	LEU	8.0
1	F	269	LEU	7.4
1	L	413	ILE	7.4
1	B	269	LEU	7.1
1	L	356	TYR	6.6
1	L	378	GLY	6.5
1	L	386	PHE	6.5
1	I	269	LEU	6.5
1	K	243	PHE	6.4

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Mol	Chain	Res	Type	RSRZ
1	L	269	LEU	6.1
1	B	265	VAL	6.0
1	L	265	VAL	5.9
1	L	243	PHE	5.9
1	B	266	THR	5.9
1	F	264	ARG	5.7
1	J	265	VAL	5.7
1	L	327	LYS	5.7
1	B	268	GLU	5.6
1	L	369	CYS	5.5
1	L	310	VAL	5.2
1	L	396	ALA	5.2
1	D	267	LEU	5.1
1	E	265	VAL	5.1
1	L	357	ILE	5.1
1	L	318	PHE	5.1
1	L	380	PHE	5.0
1	L	389	VAL	5.0
1	F	267	LEU	4.9
1	A	374	ALA	4.8
1	I	412	THR	4.7
1	C	269	LEU	4.6
1	K	48	ILE	4.6
1	E	269	LEU	4.6
1	I	265	VAL	4.6
1	A	461	ALA	4.5
1	L	332	GLY	4.5
1	A	480	TYR	4.5
1	L	422	ASN	4.5
1	I	411	LYS	4.5
1	C	267	LEU	4.5
1	B	243	PHE	4.4
1	L	465	PHE	4.4
1	A	481	GLY	4.4
1	C	270	GLY	4.4
1	G	468	TYR	4.4
1	C	265	VAL	4.4
1	L	314	ILE	4.4
1	L	400	ILE	4.3
1	L	223	GLY	4.3
1	D	32	VAL	4.3
1	L	7	ALA	4.3

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Mol	Chain	Res	Type	RSRZ
1	L	364	GLY	4.3
1	D	373	ILE	4.3
1	I	266	THR	4.2
1	L	387	GLY	4.2
1	L	377	ARG	4.1
1	A	376	ASP	4.1
1	J	245	GLY	4.1
1	L	416	VAL	4.1
1	L	477	LEU	4.1
1	D	226	PRO	4.1
1	L	244	THR	4.0
1	K	376	ASP	4.0
1	H	7	ALA	4.0
1	D	465	PHE	3.9
1	J	269	LEU	3.9
1	L	354	LEU	3.9
1	L	370	GLY	3.9
1	L	368	LEU	3.9
1	J	468	TYR	3.9
1	H	265	VAL	3.9
1	G	269	LEU	3.8
1	L	367	LEU	3.8
1	D	413	ILE	3.8
1	I	362	GLN	3.8
1	L	322	SER	3.7
1	L	362	GLN	3.7
1	L	359	THR	3.7
1	L	225	GLY	3.7
1	J	270	GLY	3.6
1	K	377	ARG	3.6
1	I	468	TYR	3.6
1	J	224	PHE	3.6
1	J	266	THR	3.6
1	K	108	LEU	3.6
1	L	285	TRP	3.6
1	A	373	ILE	3.6
1	J	225	GLY	3.6
1	L	411	LYS	3.6
1	I	268	GLU	3.5
1	D	422	ASN	3.5
1	L	268	GLU	3.5
1	C	243	PHE	3.5

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Mol	Chain	Res	Type	RSRZ
1	E	243	PHE	3.5
1	L	360	GLY	3.4
1	J	267	LEU	3.4
1	I	270	GLY	3.4
1	L	277	ILE	3.4
1	L	361	LYS	3.4
1	K	332	GLY	3.4
1	G	400	ILE	3.4
1	L	408	LEU	3.4
1	L	10	ALA	3.4
1	G	264	ARG	3.4
1	A	17	VAL	3.4
1	K	37	PHE	3.3
1	D	34	ARG	3.3
1	G	265	VAL	3.3
1	K	224	PHE	3.3
1	L	231	ALA	3.3
1	L	353	ILE	3.3
1	E	356	TYR	3.3
1	K	226	PRO	3.3
1	L	393	MET	3.3
1	L	412	THR	3.3
1	L	395	ILE	3.2
1	L	365	ALA	3.2
1	D	269	LEU	3.2
1	K	422	ASN	3.2
1	L	358	ASN	3.2
1	A	365	ALA	3.2
1	L	384	THR	3.2
1	E	267	LEU	3.2
1	L	373	ILE	3.2
1	D	461	ALA	3.2
1	L	313	ASP	3.2
1	L	376	ASP	3.2
1	E	422	ASN	3.2
1	L	338	LYS	3.1
1	F	466	GLY	3.1
1	F	265	VAL	3.1
1	A	7	ALA	3.1
1	I	408	LEU	3.1
1	D	356	TYR	3.1
1	H	468	TYR	3.1

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Mol	Chain	Res	Type	RSRZ
1	K	245	GLY	3.1
1	L	407	ILE	3.1
1	L	311	GLN	3.0
1	D	7	ALA	3.0
1	E	477	LEU	3.0
1	L	264	ARG	3.0
1	L	337	SER	3.0
1	I	243	PHE	3.0
1	B	270	GLY	3.0
1	J	243	PHE	3.0
1	H	422	ASN	2.9
1	L	385	VAL	2.9
1	D	22	ILE	2.9
1	K	400	ILE	2.9
1	D	477	LEU	2.9
1	I	389	VAL	2.9
1	G	467	GLY	2.9
1	A	386	PHE	2.9
1	K	465	PHE	2.9
1	L	419	ARG	2.9
1	D	468	TYR	2.9
1	K	193	VAL	2.8
1	L	287	VAL	2.8
1	E	264	ARG	2.8
1	L	409	LYS	2.8
1	A	422	ASN	2.8
1	D	376	ASP	2.8
1	L	215	PRO	2.8
1	I	267	LEU	2.8
1	K	401	PHE	2.8
1	A	423	SER	2.8
1	F	266	THR	2.8
1	E	468	TYR	2.8
1	D	58	ASP	2.8
1	I	422	ASN	2.8
1	D	59	VAL	2.8
1	L	188	VAL	2.8
1	L	309	PHE	2.7
1	A	314	ILE	2.7
1	I	276	ILE	2.7
1	L	276	ILE	2.7
1	D	30	ASP	2.7

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Mol	Chain	Res	Type	RSRZ
1	I	356	TYR	2.7
1	L	295	PHE	2.7
1	L	59	VAL	2.7
1	L	328	SER	2.7
1	K	57	GLU	2.7
1	L	421	ASN	2.7
1	B	238	VAL	2.7
1	J	7	ALA	2.7
1	L	237	ASP	2.7
1	I	375	ALA	2.7
1	K	52	ALA	2.7
1	A	315	TYR	2.7
1	E	7	ALA	2.6
1	E	461	ALA	2.6
1	D	423	SER	2.6
1	A	377	ARG	2.6
1	A	421	ASN	2.6
1	K	32	VAL	2.6
1	D	243	PHE	2.6
1	C	477	LEU	2.6
1	L	446	LEU	2.6
1	J	226	PRO	2.6
1	L	363	GLU	2.6
1	I	245	GLY	2.6
1	I	393	MET	2.6
1	L	72	LEU	2.6
1	I	400	ILE	2.6
1	L	238	VAL	2.6
1	A	462	GLN	2.6
1	K	244	THR	2.6
1	L	420	ALA	2.6
1	A	320	GLU	2.6
1	J	476	GLU	2.6
1	A	323	VAL	2.6
1	A	264	ARG	2.6
1	C	264	ARG	2.6
1	D	378	GLY	2.5
1	B	356	TYR	2.5
1	H	356	TYR	2.5
1	K	445	ALA	2.5
1	G	477	LEU	2.5
1	L	278	MET	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	28	TRP	2.5
1	D	61	LYS	2.5
1	L	55	ASP	2.5
1	L	317	GLU	2.5
1	L	330	VAL	2.5
1	I	271	GLY	2.5
1	D	57	GLU	2.5
1	I	316	ASP	2.5
1	L	316	ASP	2.5
1	I	318	PHE	2.5
1	D	482	LEU	2.5
1	I	264	ARG	2.5
1	J	354	LEU	2.5
1	J	467	GLY	2.5
1	K	223	GLY	2.5
1	D	29	HIS	2.5
1	A	319	VAL	2.5
1	L	8	VAL	2.5
1	L	37	PHE	2.4
1	H	267	LEU	2.4
1	I	311	GLN	2.4
1	K	47	VAL	2.4
1	A	369	CYS	2.4
1	D	463	SER	2.4
1	J	411	LYS	2.4
1	L	329	ARG	2.4
1	I	388	ASP	2.4
1	J	376	ASP	2.4
1	L	266	THR	2.4
1	K	411	LYS	2.4
1	L	226	PRO	2.4
1	H	269	LEU	2.4
1	E	268	GLU	2.4
1	B	225	GLY	2.4
1	C	496	PRO	2.4
1	I	368	LEU	2.4
1	A	417	VAL	2.4
1	A	254	GLN	2.4
1	L	447	GLN	2.4
1	G	267	LEU	2.4
1	H	243	PHE	2.4
1	A	10	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	L	189	VAL	2.3
1	L	481	GLY	2.3
1	K	72	LEU	2.3
1	B	224	PHE	2.3
1	D	389	VAL	2.3
1	K	36	THR	2.3
1	K	345	VAL	2.3
1	K	318	PHE	2.3
1	K	441	TYR	2.3
1	D	466	GLY	2.3
1	L	14	GLN	2.3
1	L	234	SER	2.3
1	L	382	GLN	2.3
1	L	17	VAL	2.3
1	F	144	ILE	2.3
1	G	263	LYS	2.3
1	K	144	ILE	2.3
1	D	224	PHE	2.3
1	G	270	GLY	2.3
1	L	315	TYR	2.3
1	A	251	ARG	2.3
1	A	392	GLY	2.2
1	A	459	PHE	2.2
1	I	312	GLU	2.2
1	I	461	ALA	2.2
1	D	464	PRO	2.2
1	L	403	PRO	2.2
1	E	466	GLY	2.2
1	K	234	SER	2.2
1	E	362	GLN	2.2
1	H	264	ARG	2.2
1	L	461	ALA	2.2
1	H	376	ASP	2.2
1	D	26	ASN	2.2
1	D	334	PRO	2.2
1	H	263	LYS	2.2
1	L	224	PHE	2.2
1	J	400	ILE	2.2
1	K	476	GLU	2.2
1	L	75	PRO	2.2
1	D	362	GLN	2.2
1	I	244	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	I	465	PHE	2.2
1	L	423	SER	2.2
1	G	356	TYR	2.2
1	H	236	GLU	2.2
1	L	281	ALA	2.2
1	H	477	LEU	2.1
1	A	279	SER	2.1
1	C	266	THR	2.1
1	L	308	THR	2.1
1	L	371	GLY	2.1
1	A	14	GLN	2.1
1	A	18	PHE	2.1
1	L	177	TRP	2.1
1	L	326	ALA	2.1
1	L	441	TYR	2.1
1	L	217	VAL	2.1
1	D	234	SER	2.1
1	F	467	GLY	2.1
1	H	362	GLN	2.1
1	A	375	ALA	2.1
1	C	7	ALA	2.1
1	D	230	ALA	2.1
1	D	478	GLY	2.1
1	F	270	GLY	2.1
1	J	46	GLU	2.1
1	K	423	SER	2.1
1	L	57	GLU	2.1
1	L	350	PHE	2.1
1	A	47	VAL	2.1
1	J	227	THR	2.1
1	D	60	ASP	2.1
1	G	146	ILE	2.1
1	A	327	LYS	2.1
1	C	467	GLY	2.1
1	K	7	ALA	2.1
1	L	230	ALA	2.1
1	D	337	SER	2.1
1	D	64	LYS	2.0
1	B	271	GLY	2.0
1	L	429	ALA	2.0
1	A	389	VAL	2.0
1	A	441	TYR	2.0

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Mol	Chain	Res	Type	RSRZ
1	G	422	ASN	2.0
1	C	400	ILE	2.0
1	I	446	LEU	2.0
1	L	463	SER	2.0
1	A	20	ASN	2.0
1	A	331	VAL	2.0
1	L	381	ILE	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 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
4	EDO	A	3116	4/4	0.64	0.33	9.30	47,49,53,54	0
4	EDO	F	3122	4/4	0.92	0.26	8.07	47,50,54,56	0
4	EDO	A	3109	4/4	0.91	0.19	5.83	47,51,51,56	0
3	GAI	I	3107	4/4	0.88	0.28	5.53	48,50,53,54	0
3	GAI	L	3108	4/4	0.56	0.37	5.44	52,53,55,57	0
4	EDO	A	3110	4/4	0.79	0.35	3.84	55,57,57,58	0
4	EDO	C	3114	4/4	0.94	0.19	3.73	43,45,46,50	0
4	EDO	C	3113	4/4	0.66	0.21	2.60	53,54,57,57	0
4	EDO	F	3120	4/4	0.95	0.22	2.49	50,52,55,55	0
3	GAI	F	3103	4/4	0.83	0.24	2.39	52,55,55,58	0
4	EDO	G	3123	4/4	0.89	0.18	2.02	51,55,55,56	0
3	GAI	A	3101	4/4	0.90	0.21	1.71	40,43,45,45	0
4	EDO	D	3126	4/4	0.85	0.17	1.63	55,58,59,59	0
4	EDO	D	3117	4/4	0.70	0.23	1.61	50,54,56,56	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
4	EDO	B	3112	4/4	0.92	0.14	1.38	38,44,48,50	0
3	GAI	F	3104	4/4	0.96	0.18	0.86	46,46,46,47	0
4	EDO	F	3121	4/4	0.96	0.13	0.85	27,42,45,46	0
3	GAI	E	3102	4/4	0.94	0.17	0.85	34,38,40,44	0
2	NA	B	3702	1/1	0.97	0.10	0.80	30,30,30,30	0
4	EDO	E	3118	4/4	0.91	0.15	0.79	39,49,51,53	0
4	EDO	J	3128	4/4	0.74	0.22	0.76	51,52,54,56	0
4	EDO	L	3129	4/4	0.87	0.14	0.55	55,55,56,57	0
3	GAI	H	3106	4/4	0.93	0.15	0.19	48,48,49,50	0
4	EDO	G	3124	4/4	0.88	0.16	0.11	41,48,51,52	0
2	NA	K	3711	1/1	0.88	0.18	0.00	73,73,73,73	0
3	GAI	G	3105	4/4	0.95	0.14	-0.09	40,45,46,47	0
2	NA	A	3701	1/1	0.77	0.14	-0.10	57,57,57,57	0
2	NA	I	3709	1/1	0.96	0.09	-0.34	36,36,36,36	0
2	NA	C	3703	1/1	0.90	0.10	-0.59	34,34,34,34	0
2	NA	D	3704	1/1	0.71	0.12	-0.82	55,55,55,55	0
2	NA	J	3710	1/1	0.75	0.10	-0.84	48,48,48,48	0
2	NA	L	3712	1/1	0.92	0.12	-0.85	60,60,60,60	0
2	NA	E	3705	1/1	0.97	0.08	-1.08	28,28,28,28	0
2	NA	H	3708	1/1	0.95	0.09	-1.90	36,36,36,36	0
2	NA	F	3706	1/1	0.97	0.07	-2.50	25,25,25,25	0
2	NA	G	3707	1/1	0.98	0.06	-3.18	37,37,37,37	0
4	EDO	G	3125	4/4	0.93	0.21	-	56,56,58,59	0
4	EDO	C	3115	4/4	0.87	0.25	-	53,53,57,59	0
4	EDO	I	3127	4/4	0.85	0.32	-	52,52,57,58	0
4	EDO	I	3119	4/4	0.83	0.25	-	53,53,57,61	0
4	EDO	A	3111	4/4	0.86	0.20	-	55,57,57,58	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.