



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:31 PM GMT

PDB ID : 1FZ4  
Title : METHANE MONOOXYGENASE HYDROXYLASE, FORM III SOAKED  
AT PH 8.5 (0.1 M TRIS)  
Authors : Whittington, D.A.; Lippard, S.J.  
Deposited on : 2000-10-03  
Resolution : 2.38 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

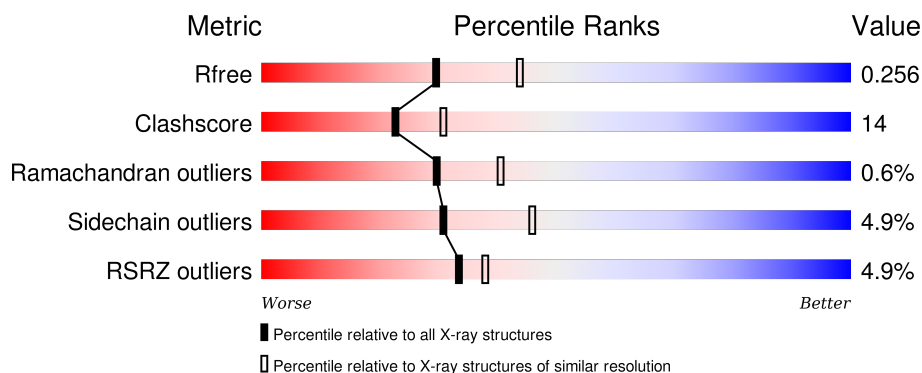
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.38 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	4019 (2.40-2.36)
Clashscore	102246	4595 (2.40-2.36)
Ramachandran outliers	100387	4520 (2.40-2.36)
Sidechain outliers	100360	4522 (2.40-2.36)
RSRZ outliers	91569	4034 (2.40-2.36)

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	527	<div> <div>5%</div> <div>69%</div> <div>24%</div> <div>• •</div> </div>
1	B	527	<div> <div>3%</div> <div>68%</div> <div>26%</div> <div>• •</div> </div>
2	C	389	<div> <div>%</div> <div>76%</div> <div>22%</div> <div>•</div> </div>
2	D	389	<div> <div>9%</div> <div>74%</div> <div>23%</div> <div>• •</div> </div>
3	E	170	<div> <div>%</div> <div>66%</div> <div>28%</div> <div>• •</div> </div>

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Mol	Chain	Length	Quality of chain
3	F	170	<div><div></div><div>14%</div><div>69%</div><div>28%</div><div>..</div></div>

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 17975 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 METHANE MONOOXYGENASE COMPONENT A, ALPHA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	511	Total	C	N	O	S	0	0	0
			4185	2677	721	769	18			
1	B	510	Total	C	N	O	S	0	0	0
			4177	2673	719	767	18			

- Molecule 2 is a protein called METHANE MONOOXYGENASE COMPONENT A, BETA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	388	Total	C	N	O	S	0	0	0
			3193	2054	551	580	8			
2	D	387	Total	C	N	O	S	0	0	0
			3183	2048	549	578	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	370	ARG	ALA	CONFLICT	UNP P18798
D	370	ARG	ALA	CONFLICT	UNP P18798

- Molecule 3 is a protein called METHANE MONOOXYGENASE COMPONENT A, GAMMA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	166	Total	C	N	O	S	0	0	0
			1368	867	246	250	5			
3	F	168	Total	C	N	O	S	0	0	0
			1386	878	250	253	5			

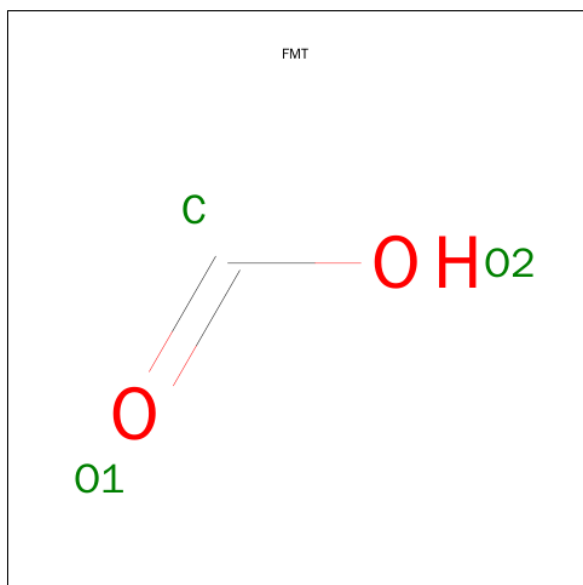
- Molecule 4 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total	Fe	0	0
			2	2		
4	A	2	Total	Fe	0	0
			2	2		

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Ca	0	0
			1	1		
5	C	2	Total	Ca	0	0
			2	2		

- Molecule 6 is FORMIC ACID (three-letter code: FMT) (formula: CH<sub>2</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			3	1	2		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	102	Total	O	0	0
			102	102		
7	B	107	Total	O	0	0
			107	107		

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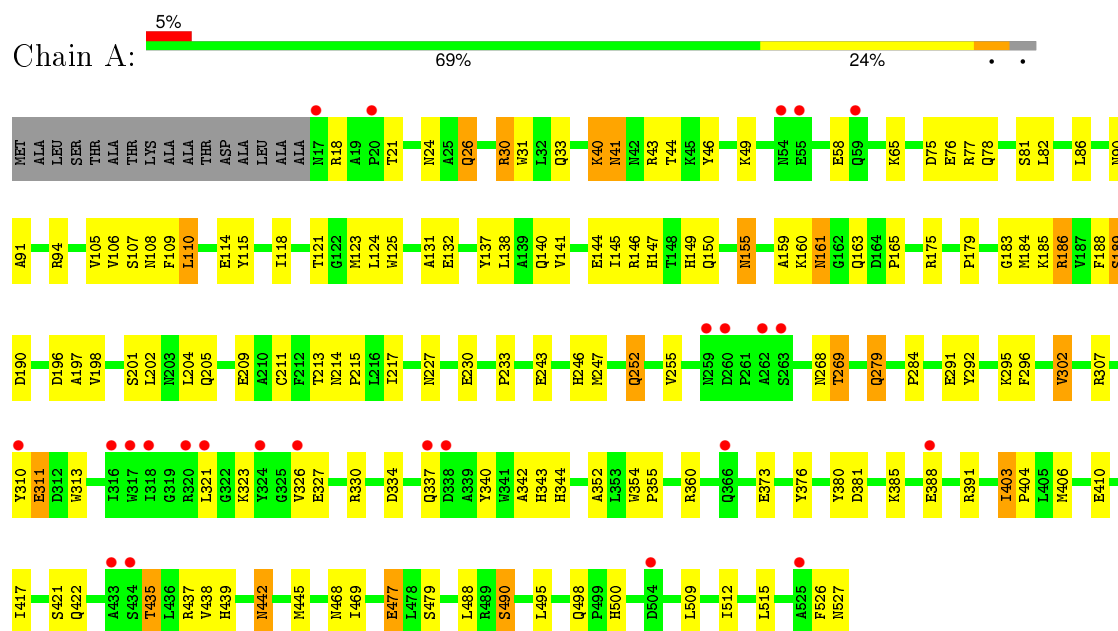
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	C	147	Total 147	O 147	0	0
7	D	41	Total 41	O 41	0	0
7	E	65	Total 65	O 65	0	0
7	F	11	Total 11	O 11	0	0

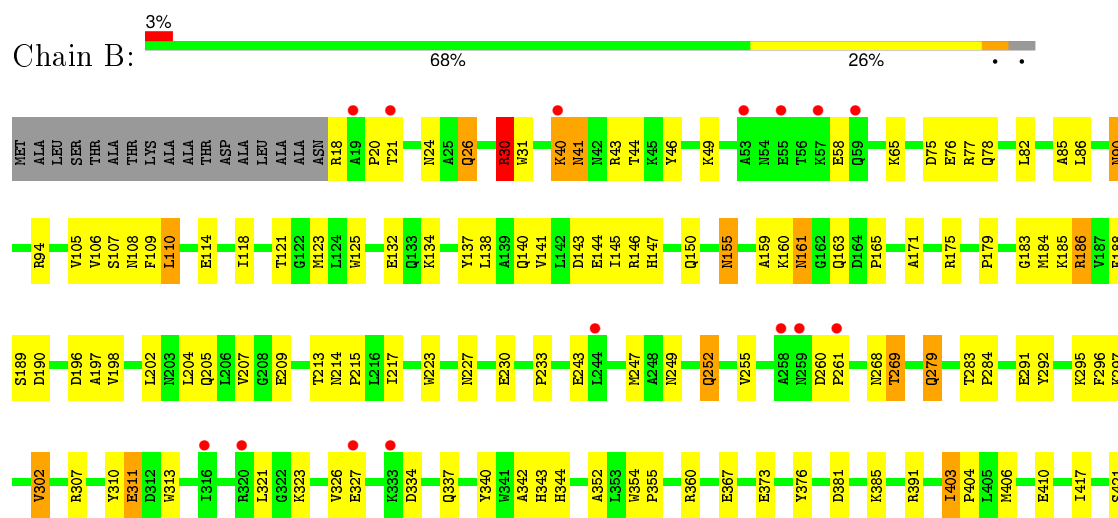
### 3 Residue-property plots

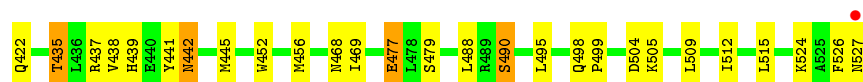
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: METHANE MONOOXYGENASE COMPONENT A, ALPHA CHAIN

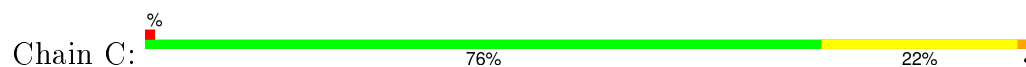


#### • Molecule 1: METHANE MONOOXYGENASE COMPONENT A, ALPHA CHAIN

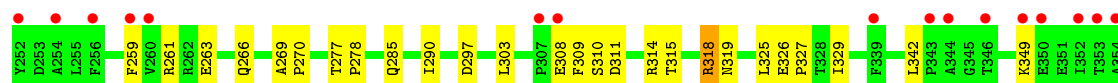
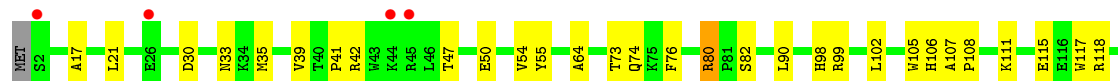
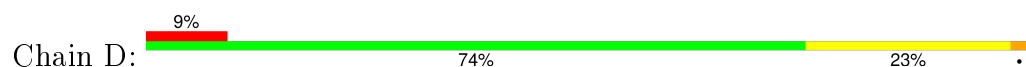




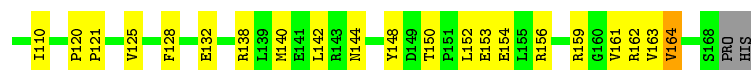
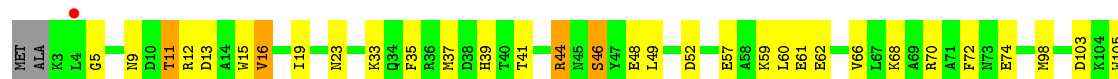
• Molecule 2: METHANE MONOOXYGENASE COMPONENT A, BETA CHAIN



• Molecule 2: METHANE MONOOXYGENASE COMPONENT A, BETA CHAIN



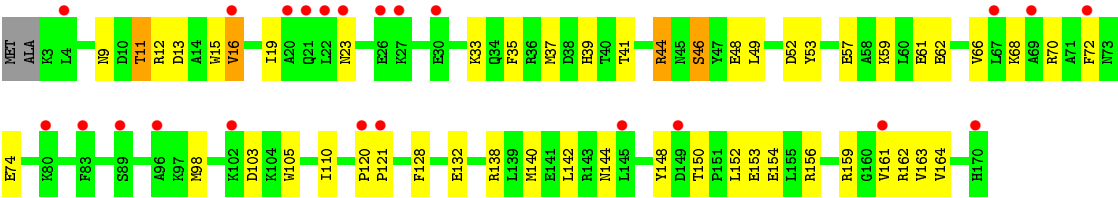
• Molecule 3: METHANE MONOOXYGENASE COMPONENT A, GAMMA CHAIN



• Molecule 3: METHANE MONOOXYGENASE COMPONENT A, GAMMA CHAIN







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	70.89Å 171.52Å 221.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.70 – 2.38 29.71 – 2.25	Depositor EDS
% Data completeness (in resolution range)	84.0 (29.70-2.38) 80.3 (29.71-2.25)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.87 (at 2.24Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.223 , 0.256 0.223 , 0.256	Depositor DCC
$R_{free}$ test set	3231 reflections (3.52%)	DCC
Wilson B-factor (Å <sup>2</sup> )	33.5	Xtriage
Anisotropy	0.275	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 41.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 117974 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	17975	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.57% 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: FMT, CA, FE

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	0/4310	0.72	10/5853 (0.2%)
1	B	0.38	0/4302	0.70	9/5842 (0.2%)
2	C	0.40	0/3289	0.58	0/4464
2	D	0.37	0/3279	0.57	0/4453
3	E	0.39	0/1396	0.60	0/1880
3	F	0.36	0/1416	0.58	0/1907
All	All	0.38	0/17992	0.65	19/24399 (0.1%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	77	ARG	NE-CZ-NH1	-15.29	112.66	120.30
1	B	30	ARG	NE-CZ-NH2	-12.81	113.89	120.30
1	A	77	ARG	NE-CZ-NH2	12.40	126.50	120.30
1	B	391	ARG	NE-CZ-NH2	-12.24	114.18	120.30
1	A	30	ARG	NE-CZ-NH1	-12.20	114.20	120.30
1	A	391	ARG	NE-CZ-NH1	-11.60	114.50	120.30
1	B	77	ARG	NE-CZ-NH2	-11.46	114.57	120.30
1	A	30	ARG	NE-CZ-NH2	10.96	125.78	120.30
1	B	391	ARG	NE-CZ-NH1	10.89	125.75	120.30
1	B	30	ARG	NE-CZ-NH1	10.62	125.61	120.30
1	A	391	ARG	NE-CZ-NH2	9.84	125.22	120.30
1	B	77	ARG	NE-CZ-NH1	9.00	124.80	120.30
1	A	30	ARG	CD-NE-CZ	6.64	132.89	123.60
1	B	30	ARG	CD-NE-CZ	6.13	132.18	123.60
1	A	279	GLN	CA-CB-CG	6.02	126.64	113.40
1	A	77	ARG	CD-NE-CZ	6.00	132.00	123.60
1	B	391	ARG	CD-NE-CZ	5.75	131.65	123.60
1	A	391	ARG	CD-NE-CZ	5.61	131.46	123.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	279	GLN	CA-CB-CG	5.17	124.78	113.40

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	4185	0	3981	125	0
1	B	4177	0	3975	133	0
2	C	3193	0	3042	87	0
2	D	3183	0	3029	87	0
3	E	1368	0	1363	49	0
3	F	1386	0	1377	48	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
5	A	1	0	0	0	0
5	C	2	0	0	0	0
6	A	3	0	1	0	0
7	A	102	0	0	4	0
7	B	107	0	0	4	0
7	C	147	0	0	4	0
7	D	41	0	0	1	0
7	E	65	0	0	0	0
7	F	11	0	0	1	0
All	All	17975	0	16768	466	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 14.

All (466) 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:B:252:GLN:HE21	1:B:252:GLN:HA	1.12	1.08
1:A:252:GLN:HE21	1:A:252:GLN:HA	1.11	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107:SER:HB3	1:B:155:ASN:HD21	1.16	1.06
2:C:270:PRO:HB3	2:D:270:PRO:HB3	1.39	1.04
1:A:107:SER:HB3	1:A:155:ASN:HD21	1.20	1.04
2:D:146:ASN:HD21	2:D:197:ARG:HH21	1.00	0.97
2:C:146:ASN:HD21	2:C:197:ARG:HH21	0.97	0.96
1:B:78:GLN:HE22	1:B:150:GLN:HE21	1.10	0.93
1:A:44:THR:HG22	1:A:46:TYR:H	1.32	0.93
2:C:319:ASN:HA	3:E:74:GLU:OE1	1.70	0.91
1:A:78:GLN:HE22	1:A:150:GLN:HE21	1.03	0.91
1:B:160:LYS:C	1:B:161:ASN:HD22	1.74	0.91
1:B:44:THR:HG22	1:B:46:TYR:H	1.32	0.90
1:A:209:GLU:HA	1:A:213:THR:HB	1.52	0.90
2:C:146:ASN:ND2	2:C:197:ARG:HH21	1.69	0.89
2:C:261:ARG:HE	2:C:285:GLN:HE22	1.18	0.88
2:D:146:ASN:ND2	2:D:197:ARG:HH21	1.72	0.87
1:A:78:GLN:NE2	1:A:150:GLN:HE21	1.72	0.87
1:B:352:ALA:HA	1:B:404:PRO:HB2	1.57	0.87
2:D:261:ARG:HE	2:D:285:GLN:HE22	1.19	0.86
1:A:352:ALA:HA	1:A:404:PRO:HB2	1.55	0.86
1:A:41:ASN:HD22	1:A:41:ASN:N	1.73	0.86
1:B:107:SER:CB	1:B:155:ASN:HD21	1.89	0.84
2:D:99:ARG:HH11	2:D:99:ARG:HG2	1.41	0.83
1:B:252:GLN:NE2	1:B:252:GLN:HA	1.92	0.83
1:A:252:GLN:HA	1:A:252:GLN:NE2	1.93	0.82
1:A:108:ASN:HD21	1:A:175:ARG:HH11	1.27	0.81
1:B:41:ASN:HD22	1:B:41:ASN:N	1.78	0.80
1:A:160:LYS:C	1:A:161:ASN:HD22	1.85	0.79
1:B:78:GLN:NE2	1:B:150:GLN:HE21	1.81	0.79
2:D:319:ASN:HA	3:F:74:GLU:OE1	1.84	0.78
2:C:102:LEU:CD1	2:C:290:ILE:HG23	2.15	0.77
1:B:108:ASN:HD21	1:B:175:ARG:HH11	1.32	0.77
2:C:146:ASN:HD21	2:C:197:ARG:NH2	1.81	0.77
1:B:107:SER:HB3	1:B:155:ASN:ND2	1.97	0.76
1:A:107:SER:CB	1:A:155:ASN:HD21	1.96	0.76
2:D:102:LEU:CD1	2:D:290:ILE:HG23	2.17	0.74
2:C:315:THR:HA	2:C:318:ARG:NH1	2.02	0.74
1:B:209:GLU:HA	1:B:213:THR:HB	1.68	0.74
2:C:315:THR:HA	2:C:318:ARG:HH12	1.53	0.74
1:A:355:PRO:HG2	1:A:403:ILE:HD11	1.70	0.73
3:F:98:MET:HE2	3:F:110:ILE:HB	1.70	0.73
1:A:78:GLN:HE22	1:A:150:GLN:NE2	1.84	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:98:MET:HE2	3:E:110:ILE:HB	1.70	0.73
2:D:377:ARG:HG2	7:D:422:HOH:O	1.88	0.73
2:D:315:THR:HA	2:D:318:ARG:NH1	2.03	0.72
1:B:355:PRO:HG2	1:B:403:ILE:HD11	1.69	0.72
2:D:261:ARG:HE	2:D:285:GLN:NE2	1.88	0.72
2:C:261:ARG:HE	2:C:285:GLN:NE2	1.88	0.71
1:A:155:ASN:HD22	1:A:155:ASN:N	1.87	0.71
2:D:315:THR:HA	2:D:318:ARG:HH12	1.55	0.70
1:B:75:ASP:OD2	1:B:146:ARG:NH1	2.25	0.69
1:B:161:ASN:HD22	1:B:161:ASN:N	1.86	0.69
3:F:41:THR:O	3:F:44:ARG:HD2	1.93	0.69
2:C:99:ARG:HH11	2:C:99:ARG:HG2	1.58	0.69
1:B:227:ASN:HD21	1:B:295:LYS:H	1.39	0.69
1:A:252:GLN:HE21	1:A:252:GLN:CA	1.98	0.69
3:F:57:GLU:O	3:F:61:GLU:HG3	1.93	0.68
1:A:40:LYS:HB3	1:A:41:ASN:HD22	1.59	0.67
2:C:146:ASN:HD22	2:C:197:ARG:HE	1.41	0.67
3:F:68:LYS:HG3	3:F:72:PHE:CD2	2.29	0.67
1:B:49:LYS:CE	3:F:144:ASN:HD22	2.08	0.67
3:F:9:ASN:OD1	3:F:11:THR:HG22	1.95	0.67
1:A:118:ILE:HD13	1:A:145:ILE:HG12	1.75	0.66
1:B:49:LYS:HE3	3:F:144:ASN:HD22	1.61	0.66
2:D:76:PHE:HZ	2:D:168:ARG:HH12	1.41	0.66
2:C:376:ASP:HB3	2:C:379:GLN:HG3	1.78	0.66
1:B:179:PRO:HB3	1:B:469:ILE:HD13	1.76	0.66
3:E:57:GLU:O	3:E:61:GLU:HG3	1.97	0.65
2:D:376:ASP:HB3	2:D:379:GLN:HG3	1.78	0.65
1:A:107:SER:HB3	1:A:155:ASN:ND2	2.02	0.65
1:B:155:ASN:HD22	1:B:155:ASN:N	1.93	0.65
2:D:99:ARG:HG2	2:D:99:ARG:NH1	2.10	0.65
1:A:227:ASN:HD21	1:A:295:LYS:H	1.44	0.65
2:D:107:ALA:HB3	2:D:108:PRO:HD3	1.77	0.65
1:B:185:LYS:O	1:B:189:SER:HB2	1.97	0.64
1:B:360:ARG:HG2	1:B:498:GLN:HB2	1.78	0.64
3:F:39:HIS:CD2	3:F:49:LEU:HD12	2.32	0.64
1:B:269:THR:HG22	3:F:148:TYR:CE1	2.32	0.64
1:B:165:PRO:HG3	7:B:5010:HOH:O	1.96	0.64
3:E:138:ARG:HH21	3:E:142:LEU:HD21	1.62	0.63
1:B:268:ASN:HD21	1:B:327:GLU:H	1.46	0.63
1:A:292:TYR:OH	1:A:344:HIS:HD2	1.81	0.63
2:C:326:GLU:HB3	2:C:327:PRO:HD3	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:31:TRP:CH2	2:D:210:SER:HA	2.34	0.63
2:C:107:ALA:HB3	2:C:108:PRO:HD3	1.79	0.63
3:E:68:LYS:HG3	3:E:72:PHE:CD2	2.33	0.63
2:C:47:THR:OG1	2:C:50:GLU:HG3	1.98	0.63
3:E:41:THR:O	3:E:44:ARG:HD2	1.97	0.63
2:D:47:THR:OG1	2:D:50:GLU:HG3	1.99	0.63
1:B:160:LYS:HG2	1:B:161:ASN:ND2	2.14	0.63
1:A:185:LYS:O	1:A:189:SER:HB2	1.99	0.62
1:A:108:ASN:ND2	1:A:175:ARG:HH11	1.98	0.62
2:D:146:ASN:HD22	2:D:197:ARG:HE	1.48	0.62
2:D:111:LYS:O	2:D:115:GLU:HG3	1.98	0.62
1:A:268:ASN:HD21	1:A:327:GLU:H	1.47	0.62
1:A:495:LEU:HD11	1:A:512:ILE:HG13	1.80	0.62
1:B:495:LEU:HD11	1:B:512:ILE:HG13	1.81	0.62
1:B:292:TYR:OH	1:B:344:HIS:HD2	1.83	0.61
1:B:41:ASN:ND2	1:B:41:ASN:N	2.48	0.61
1:B:439:HIS:HD2	3:F:163:VAL:HA	1.66	0.61
2:C:76:PHE:HZ	2:C:168:ARG:HH12	1.48	0.61
1:A:355:PRO:HG2	1:A:403:ILE:CD1	2.31	0.61
2:D:195:LEU:O	2:D:195:LEU:HD23	2.00	0.61
1:A:477:GLU:OE2	1:A:479:SER:N	2.34	0.60
2:D:146:ASN:HD21	2:D:197:ARG:NH2	1.85	0.60
2:C:111:LYS:O	2:C:115:GLU:HG3	2.00	0.60
1:B:161:ASN:N	1:B:161:ASN:ND2	2.49	0.60
2:D:326:GLU:HB3	2:D:327:PRO:HD3	1.84	0.60
1:A:360:ARG:HG2	1:A:498:GLN:HB2	1.84	0.60
3:E:39:HIS:CD2	3:E:49:LEU:HD12	2.36	0.60
1:B:252:GLN:HE21	1:B:252:GLN:CA	2.00	0.59
1:A:214:ASN:HB3	1:A:215:PRO:HD3	1.84	0.59
1:B:138:LEU:HD22	2:D:160:PHE:CZ	2.37	0.59
1:A:302:VAL:HG13	1:A:376:TYR:HE2	1.66	0.59
3:E:9:ASN:OD1	3:E:11:THR:HG22	2.03	0.59
3:F:41:THR:HG22	7:F:176:HOH:O	2.03	0.59
1:A:438:VAL:HG12	3:E:164:VAL:CG2	2.32	0.59
2:C:223:VAL:HG23	7:C:5064:HOH:O	2.01	0.59
1:A:179:PRO:HB3	1:A:469:ILE:HD13	1.85	0.59
2:D:187:ILE:O	2:D:191:GLN:HG3	2.01	0.59
1:A:40:LYS:HB3	1:A:41:ASN:ND2	2.18	0.59
1:B:477:GLU:OE2	1:B:479:SER:N	2.36	0.59
1:B:214:ASN:HB3	1:B:215:PRO:HD3	1.85	0.59
2:D:139:THR:O	2:D:143:GLU:HB3	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:195:LEU:O	2:C:195:LEU:HD23	2.03	0.58
2:C:139:THR:O	2:C:143:GLU:HB3	2.03	0.58
1:B:227:ASN:HD21	1:B:296:PHE:H	1.51	0.58
2:C:376:ASP:CB	2:C:379:GLN:HG3	2.33	0.58
1:A:161:ASN:N	1:A:161:ASN:HD22	1.98	0.57
1:B:355:PRO:HG2	1:B:403:ILE:CD1	2.33	0.57
1:A:160:LYS:HA	2:C:33:ASN:HB2	1.85	0.57
1:A:435:THR:CG2	1:A:437:ARG:HE	2.17	0.57
1:B:302:VAL:HG13	1:B:376:TYR:HE2	1.68	0.57
1:B:49:LYS:HD3	3:F:140:MET:HB3	1.87	0.57
2:D:349:LYS:HE3	2:D:385:LEU:HD12	1.86	0.57
2:C:308:GLU:HG2	2:C:309:PHE:CE1	2.40	0.57
1:B:30:ARG:O	1:B:30:ARG:HD3	2.05	0.57
2:C:349:LYS:HE3	2:C:385:LEU:CD1	2.34	0.57
1:A:41:ASN:N	1:A:41:ASN:ND2	2.46	0.57
2:C:99:ARG:HG2	2:C:99:ARG:NH1	2.19	0.57
1:B:438:VAL:HG12	3:F:164:VAL:CG2	2.35	0.57
1:B:435:THR:CG2	1:B:437:ARG:HE	2.18	0.56
1:A:490:SER:OG	2:C:30:ASP:OD1	2.23	0.56
1:B:78:GLN:HE22	1:B:150:GLN:NE2	1.91	0.56
2:C:102:LEU:HD13	2:C:290:ILE:HG23	1.85	0.56
3:F:138:ARG:HH21	3:F:142:LEU:HD21	1.70	0.56
2:D:376:ASP:CB	2:D:379:GLN:HG3	2.36	0.56
2:C:105:TRP:O	2:C:108:PRO:HD2	2.06	0.56
1:B:526:PHE:O	1:B:527:ASN:ND2	2.36	0.56
3:F:15:TRP:O	3:F:19:ILE:HG13	2.05	0.56
1:A:123:MET:CE	1:A:197:ALA:HA	2.35	0.56
1:A:160:LYS:HG2	1:A:161:ASN:ND2	2.21	0.56
2:C:356:LEU:HD21	2:C:384:VAL:HG11	1.88	0.56
3:F:13:ASP:HA	3:F:16:VAL:HG23	1.88	0.56
1:A:373:GLU:OE1	1:A:373:GLU:HA	2.05	0.56
3:F:41:THR:O	3:F:44:ARG:CD	2.54	0.56
1:A:439:HIS:HD2	3:E:163:VAL:HA	1.70	0.56
1:A:58:GLU:HB3	1:A:132:GLU:O	2.05	0.56
2:D:98:HIS:HD2	2:D:297:ASP:OD1	1.89	0.56
1:B:108:ASN:ND2	1:B:175:ARG:HH11	2.02	0.55
1:B:184:MET:HE1	1:B:188:PHE:CD2	2.41	0.55
1:A:49:LYS:CE	3:E:144:ASN:HD22	2.18	0.55
1:B:442:ASN:HD22	1:B:442:ASN:C	2.08	0.55
1:B:373:GLU:HA	1:B:373:GLU:OE1	2.05	0.55
1:B:140:GLN:O	1:B:144:GLU:HG2	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:35:PHE:CE1	3:E:39:HIS:HD2	2.24	0.55
2:D:349:LYS:HE3	2:D:385:LEU:CD1	2.36	0.55
3:E:41:THR:O	3:E:44:ARG:CD	2.55	0.55
1:A:75:ASP:OD2	1:A:146:ARG:NH1	2.40	0.55
1:B:227:ASN:ND2	1:B:295:LYS:H	2.04	0.54
3:F:46:SER:OG	3:F:48:GLU:HG2	2.08	0.54
1:A:91:ALA:HB1	7:A:6094:HOH:O	2.07	0.54
2:C:376:ASP:CG	2:C:379:GLN:HG3	2.27	0.54
1:A:438:VAL:HG12	3:E:164:VAL:HG21	1.90	0.54
1:A:76:GLU:OE2	1:B:76:GLU:HG2	2.07	0.54
1:B:146:ARG:HB2	2:D:106:HIS:CE1	2.43	0.54
3:E:161:VAL:HG23	3:E:162:ARG:N	2.23	0.54
3:F:33:LYS:O	3:F:37:MET:HG2	2.08	0.54
1:A:114:GLU:CD	1:A:147:HIS:HB3	2.27	0.54
1:A:49:LYS:HE3	3:E:144:ASN:HD22	1.72	0.54
1:A:186:ARG:HA	2:C:73:THR:OG1	2.08	0.54
2:D:105:TRP:O	2:D:108:PRO:HD2	2.08	0.53
1:A:123:MET:HE1	1:A:197:ALA:HA	1.89	0.53
2:C:349:LYS:HE3	2:C:385:LEU:HD12	1.91	0.53
1:A:196:ASP:HB2	3:E:140:MET:SD	2.47	0.53
2:C:146:ASN:ND2	2:C:197:ARG:NH2	2.48	0.53
2:D:80:ARG:HB2	3:F:132:GLU:OE1	2.08	0.53
1:A:161:ASN:N	1:A:161:ASN:ND2	2.56	0.53
2:D:153:LEU:C	2:D:153:LEU:HD12	2.29	0.53
1:A:500:HIS:HB2	7:A:6060:HOH:O	2.09	0.53
2:D:90:LEU:HD13	2:D:303:LEU:HD13	1.91	0.53
1:B:160:LYS:HA	2:D:33:ASN:HB2	1.90	0.53
2:D:102:LEU:HD13	2:D:290:ILE:HG23	1.88	0.53
3:E:12:ARG:O	3:E:16:VAL:HG23	2.08	0.53
2:D:308:GLU:HG2	2:D:309:PHE:CE1	2.43	0.53
1:B:445:MET:HE1	1:B:526:PHE:O	2.09	0.52
2:C:187:ILE:O	2:C:191:GLN:HG3	2.09	0.52
3:F:159:ARG:HG3	3:F:161:VAL:HG12	1.91	0.52
3:E:15:TRP:O	3:E:19:ILE:HG13	2.08	0.52
2:C:98:HIS:HD2	2:C:297:ASP:OD1	1.92	0.52
1:A:442:ASN:HD22	1:A:442:ASN:C	2.11	0.52
1:A:284:PRO:HB3	1:A:342:ALA:HB1	1.91	0.52
1:B:490:SER:OG	2:D:30:ASP:OD1	2.27	0.52
7:B:5016:HOH:O	3:F:46:SER:HA	2.09	0.52
3:E:13:ASP:HA	3:E:16:VAL:HG23	1.92	0.52
1:A:209:GLU:HA	1:A:213:THR:CB	2.34	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:105:VAL:O	1:B:109:PHE:HB2	2.10	0.52
2:C:54:VAL:O	2:C:55:TYR:HB2	2.09	0.52
3:E:159:ARG:HG3	3:E:161:VAL:HG12	1.91	0.51
2:C:310:SER:O	2:C:314:ARG:HG3	2.10	0.51
7:C:5017:HOH:O	3:E:125:VAL:HG22	2.10	0.51
1:B:291:GLU:OE1	1:B:343:HIS:HE1	1.94	0.51
1:A:184:MET:HE1	1:A:188:PHE:CD2	2.45	0.51
3:E:33:LYS:O	3:E:37:MET:HG2	2.10	0.51
2:C:309:PHE:CZ	3:E:62:GLU:HG2	2.45	0.51
1:A:190:ASP:HB3	2:C:74:GLN:O	2.10	0.51
1:A:381:ASP:HA	1:A:385:LYS:HE2	1.91	0.51
1:A:155:ASN:ND2	1:A:155:ASN:N	2.57	0.51
2:D:82:SER:O	2:D:168:ARG:NH2	2.44	0.51
1:A:435:THR:HG23	1:A:437:ARG:HE	1.74	0.51
3:F:161:VAL:HG23	3:F:162:ARG:N	2.25	0.51
3:E:46:SER:OG	3:E:48:GLU:HG2	2.11	0.51
1:B:143:ASP:O	1:B:146:ARG:HB3	2.11	0.51
2:C:213:VAL:HG23	7:C:5068:HOH:O	2.11	0.50
1:B:252:GLN:HE22	1:B:255:VAL:HG21	1.75	0.50
2:C:365:GLU:HA	2:C:365:GLU:OE1	2.11	0.50
1:B:159:ALA:O	2:D:33:ASN:HB2	2.11	0.50
2:C:102:LEU:HD12	2:C:290:ILE:HG23	1.92	0.50
3:F:61:GLU:HB3	3:F:121:PRO:HD3	1.94	0.50
1:A:227:ASN:HD21	1:A:296:PHE:H	1.60	0.50
1:B:381:ASP:HA	1:B:385:LYS:HE2	1.93	0.50
1:B:107:SER:CB	1:B:155:ASN:ND2	2.68	0.50
1:B:160:LYS:HE3	1:B:161:ASN:HD21	1.77	0.50
1:A:140:GLN:O	1:A:144:GLU:HG2	2.12	0.50
2:C:365:GLU:CA	2:C:365:GLU:OE1	2.60	0.50
1:A:163:GLN:HG2	7:A:6095:HOH:O	2.12	0.50
1:A:31:TRP:CH2	2:C:210:SER:HA	2.46	0.50
1:B:160:LYS:HG2	1:B:161:ASN:HD21	1.77	0.49
1:A:439:HIS:CG	3:E:161:VAL:HG21	2.47	0.49
2:C:90:LEU:HD13	2:C:303:LEU:HD13	1.93	0.49
1:A:380:TYR:HE2	1:A:388:GLU:OE2	1.95	0.49
2:D:161:ASN:HB3	2:D:235:TRP:CE2	2.47	0.49
2:D:143:GLU:O	2:D:147:ARG:HB3	2.12	0.49
1:A:24:ASN:OD1	1:A:26:GLN:HG3	2.12	0.49
2:C:161:ASN:HB3	2:C:235:TRP:CE2	2.47	0.49
2:D:310:SER:O	2:D:314:ARG:HG3	2.12	0.49
1:B:24:ASN:OD1	1:B:26:GLN:HG3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:228:ARG:O	2:D:232:GLU:HG3	2.12	0.49
1:A:138:LEU:HD22	2:C:160:PHE:CZ	2.48	0.49
3:F:152:LEU:O	3:F:156:ARG:HG3	2.12	0.49
2:D:54:VAL:O	2:D:55:TYR:HB2	2.12	0.49
1:A:227:ASN:ND2	1:A:295:LYS:H	2.11	0.49
1:B:307:ARG:HA	1:B:311:GLU:HG3	1.94	0.49
1:A:302:VAL:HG11	1:A:340:TYR:CE1	2.47	0.49
2:C:153:LEU:C	2:C:153:LEU:HD12	2.33	0.49
1:B:40:LYS:HB3	1:B:41:ASN:ND2	2.28	0.49
2:D:385:LEU:C	2:D:387:GLY:H	2.15	0.49
1:A:213:THR:O	1:A:217:ILE:HG12	2.13	0.49
2:C:143:GLU:O	2:C:147:ARG:HB3	2.12	0.49
1:B:144:GLU:HA	1:B:144:GLU:OE2	2.13	0.49
1:B:284:PRO:HB3	1:B:342:ALA:HB1	1.94	0.49
1:A:137:TYR:O	1:A:141:VAL:HG23	2.13	0.49
1:A:488:LEU:HD21	1:A:509:LEU:HD13	1.95	0.49
1:A:417:ILE:HG13	1:A:468:ASN:HB2	1.95	0.49
1:A:230:GLU:C	1:A:233:PRO:HD2	2.34	0.49
1:B:367:GLU:HB2	7:B:5082:HOH:O	2.12	0.48
3:E:61:GLU:HB3	3:E:121:PRO:HD3	1.95	0.48
1:B:302:VAL:HG11	1:B:340:TYR:CE1	2.48	0.48
1:A:146:ARG:HB2	2:C:106:HIS:CE1	2.48	0.48
2:C:80:ARG:HB2	3:E:132:GLU:OE1	2.13	0.48
3:E:150:THR:HG23	3:E:154:GLU:HG2	1.94	0.48
1:A:252:GLN:HE22	1:A:255:VAL:HG21	1.78	0.48
1:A:269:THR:HG22	3:E:148:TYR:CE1	2.48	0.48
1:B:435:THR:HG23	1:B:437:ARG:HE	1.78	0.48
2:C:201:ALA:HA	2:C:207:PHE:HB3	1.95	0.48
1:B:30:ARG:HD3	1:B:30:ARG:C	2.33	0.48
1:A:307:ARG:HA	1:A:311:GLU:HG3	1.94	0.48
1:B:196:ASP:HB2	3:F:140:MET:SD	2.54	0.48
1:A:124:LEU:HD21	1:A:201:SER:HB2	1.96	0.48
1:B:118:ILE:HD13	1:B:145:ILE:HG12	1.96	0.48
1:B:141:VAL:O	1:B:145:ILE:HG13	2.14	0.48
2:D:146:ASN:ND2	2:D:197:ARG:NH2	2.52	0.48
3:E:66:VAL:HG12	3:E:70:ARG:HH21	1.79	0.48
1:B:441:TYR:HE2	3:F:159:ARG:O	1.97	0.47
1:B:198:VAL:O	1:B:202:LEU:HG	2.14	0.47
3:F:98:MET:CE	3:F:110:ILE:HB	2.42	0.47
1:B:165:PRO:HD2	2:D:30:ASP:HB3	1.96	0.47
1:A:49:LYS:HD3	3:E:140:MET:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:269:ALA:HB3	2:C:270:PRO:HD3	1.95	0.47
3:E:103:ASP:OD1	3:E:105:TRP:HB2	2.14	0.47
1:B:417:ILE:HG13	1:B:468:ASN:HB2	1.96	0.47
2:C:82:SER:O	2:C:168:ARG:NH2	2.46	0.47
1:A:110:LEU:O	1:A:114:GLU:HG2	2.14	0.47
3:F:66:VAL:HG12	3:F:70:ARG:HH21	1.80	0.47
2:D:263:GLU:HA	2:D:263:GLU:OE2	2.15	0.47
3:F:11:THR:CG2	3:F:12:ARG:N	2.78	0.47
1:B:20:PRO:HG3	2:D:129:ALA:HB2	1.96	0.47
2:D:201:ALA:HA	2:D:207:PHE:HB3	1.97	0.47
2:C:277:THR:HB	2:C:278:PRO:HD3	1.96	0.47
2:D:156:GLU:HA	2:D:156:GLU:OE2	2.14	0.47
3:F:35:PHE:CE1	3:F:39:HIS:HD2	2.33	0.47
1:B:524:LYS:O	1:B:527:ASN:HB2	2.14	0.47
1:A:165:PRO:HD2	2:C:30:ASP:HB3	1.96	0.47
2:D:235:TRP:CD1	2:D:235:TRP:C	2.88	0.47
2:C:112:ASP:OD1	2:D:118:ARG:NH2	2.48	0.47
3:F:153:GLU:CD	3:F:153:GLU:H	2.16	0.47
2:D:39:VAL:O	2:D:41:PRO:HD3	2.15	0.47
1:A:65:LYS:HB3	2:C:117:TRP:CG	2.50	0.47
1:A:243:GLU:O	1:A:247:MET:HG2	2.15	0.47
1:B:406:MET:O	1:B:410:GLU:HG3	2.15	0.47
1:A:438:VAL:HB	3:E:164:VAL:HG23	1.96	0.47
3:E:152:LEU:O	3:E:156:ARG:HG3	2.15	0.47
1:A:183:GLY:HA2	1:A:422:GLN:HB2	1.97	0.47
1:B:230:GLU:C	1:B:233:PRO:HD2	2.35	0.47
3:F:13:ASP:HA	3:F:16:VAL:CG2	2.45	0.46
1:B:137:TYR:O	1:B:141:VAL:HG23	2.15	0.46
2:C:326:GLU:CB	2:C:327:PRO:HD3	2.46	0.46
1:B:65:LYS:HB3	2:D:117:TRP:CG	2.50	0.46
2:C:193:ILE:HA	7:C:5058:HOH:O	2.15	0.46
7:A:6004:HOH:O	2:C:70:GLY:HA3	2.14	0.46
1:B:171:ALA:O	1:B:175:ARG:HD3	2.16	0.46
2:C:235:TRP:CD1	2:C:235:TRP:C	2.88	0.46
2:C:42:ARG:HB2	2:C:99:ARG:HG3	1.97	0.46
1:B:183:GLY:HA2	1:B:422:GLN:HB2	1.98	0.46
2:D:376:ASP:CG	2:D:379:GLN:HG3	2.36	0.46
1:A:211:CYS:HB2	1:A:313:TRP:CD1	2.50	0.46
3:F:159:ARG:CG	3:F:161:VAL:HG12	2.46	0.46
1:B:134:LYS:HD3	2:D:161:ASN:HD21	1.81	0.46
3:F:120:PRO:HD3	3:F:128:PHE:CG	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:144:PHE:CZ	2:D:342:LEU:HD23	2.51	0.46
1:B:488:LEU:HD21	1:B:509:LEU:HD13	1.98	0.46
3:F:103:ASP:OD1	3:F:105:TRP:HB2	2.16	0.46
1:A:403:ILE:HD12	1:A:515:LEU:CD1	2.46	0.45
2:D:326:GLU:CB	2:D:327:PRO:HD3	2.46	0.45
1:A:33:GLN:HA	1:A:131:ALA:HB3	1.99	0.45
1:A:321:LEU:O	1:A:326:VAL:HB	2.15	0.45
2:D:385:LEU:C	2:D:387:GLY:N	2.69	0.45
1:B:108:ASN:HD21	1:B:175:ARG:CD	2.30	0.45
2:D:102:LEU:HD12	2:D:290:ILE:HG23	1.94	0.45
1:B:354:TRP:CG	1:B:355:PRO:HD3	2.52	0.45
3:E:12:ARG:O	3:E:16:VAL:CG2	2.64	0.45
1:A:105:VAL:O	1:A:109:PHE:HB2	2.17	0.45
2:D:259:PHE:CE1	2:D:356:LEU:HD12	2.51	0.45
2:C:146:ASN:ND2	2:C:197:ARG:HE	2.13	0.45
2:D:356:LEU:HD21	2:D:384:VAL:HG11	1.99	0.45
1:B:155:ASN:ND2	1:B:155:ASN:N	2.61	0.45
1:A:186:ARG:HD3	1:A:186:ARG:C	2.36	0.45
1:A:291:GLU:OE1	1:A:343:HIS:HE1	2.00	0.45
1:A:354:TRP:CG	1:A:355:PRO:HD3	2.52	0.45
1:A:198:VAL:O	1:A:202:LEU:HG	2.17	0.45
2:C:39:VAL:O	2:C:41:PRO:HD3	2.16	0.45
3:F:11:THR:HG21	3:F:52:ASP:OD2	2.18	0.44
2:C:362:ASP:O	2:C:365:GLU:HB2	2.17	0.44
3:E:153:GLU:H	3:E:153:GLU:CD	2.20	0.44
3:F:12:ARG:O	3:F:16:VAL:HG23	2.16	0.44
1:A:160:LYS:HE3	1:A:161:ASN:HD21	1.80	0.44
1:B:495:LEU:HD11	1:B:512:ILE:CG1	2.47	0.44
1:B:213:THR:O	1:B:217:ILE:HG12	2.18	0.44
3:F:98:MET:HG3	3:F:138:ARG:HG2	2.00	0.44
1:B:123:MET:CE	1:B:197:ALA:HA	2.47	0.44
1:B:321:LEU:O	1:B:326:VAL:HB	2.17	0.44
1:A:445:MET:HE1	1:A:526:PHE:O	2.17	0.44
3:E:98:MET:HG3	3:E:138:ARG:HG2	2.00	0.44
1:A:477:GLU:OE2	1:A:477:GLU:C	2.56	0.44
1:A:18:ARG:HD2	2:C:131:GLY:HA3	1.99	0.44
3:E:159:ARG:CG	3:E:161:VAL:HG12	2.48	0.44
1:A:106:VAL:O	1:A:110:LEU:HB2	2.18	0.44
1:A:115:TYR:OH	2:C:173:ASP:HA	2.17	0.44
1:A:81:SER:HB3	1:B:85:ALA:HB2	2.00	0.44
3:F:150:THR:HG23	3:F:154:GLU:HG2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:98:HIS:CD2	2:D:297:ASP:OD1	2.70	0.44
1:A:114:GLU:O	1:A:144:GLU:HB3	2.18	0.44
3:E:11:THR:CG2	3:E:12:ARG:N	2.81	0.43
1:B:186:ARG:HD3	1:B:186:ARG:C	2.38	0.43
1:A:76:GLU:HG2	1:B:76:GLU:OE2	2.18	0.43
1:B:106:VAL:O	1:B:110:LEU:HB2	2.18	0.43
3:F:12:ARG:O	3:F:16:VAL:CG2	2.66	0.43
1:B:243:GLU:O	1:B:247:MET:HG2	2.18	0.43
1:A:406:MET:O	1:A:410:GLU:HG3	2.18	0.43
1:A:108:ASN:HD21	1:A:175:ARG:CD	2.32	0.43
1:B:438:VAL:HG12	3:F:164:VAL:HG21	2.00	0.43
2:D:195:LEU:C	2:D:195:LEU:HD23	2.38	0.43
2:C:144:PHE:CZ	2:C:342:LEU:HD23	2.53	0.43
1:B:190:ASP:HB3	2:D:74:GLN:O	2.18	0.43
2:D:269:ALA:HB3	2:D:270:PRO:HD3	2.00	0.43
2:D:365:GLU:OE1	2:D:365:GLU:HA	2.18	0.43
3:F:13:ASP:CA	3:F:16:VAL:HG23	2.48	0.43
3:E:19:ILE:HG12	3:E:60:LEU:HD13	2.01	0.43
2:D:146:ASN:ND2	2:D:197:ARG:HE	2.16	0.43
1:A:204:LEU:HG	1:A:205:GLN:HG3	2.01	0.43
1:B:40:LYS:HB3	1:B:41:ASN:HD22	1.81	0.43
2:C:270:PRO:CB	2:D:270:PRO:HB3	2.28	0.43
1:B:223:TRP:CZ3	1:B:297:LYS:HA	2.54	0.43
1:A:495:LEU:HD11	1:A:512:ILE:CG1	2.49	0.42
2:D:309:PHE:CZ	3:F:62:GLU:HG2	2.54	0.42
2:C:98:HIS:CD2	2:C:297:ASP:OD1	2.72	0.42
1:B:421:SER:O	1:B:422:GLN:HB2	2.19	0.42
1:B:186:ARG:HA	2:D:73:THR:OG1	2.19	0.42
2:C:228:ARG:O	2:C:232:GLU:HG3	2.19	0.42
2:D:325:LEU:O	2:D:329:ILE:HG13	2.19	0.42
3:E:59:LYS:HA	3:E:59:LYS:HD3	1.86	0.42
1:A:146:ARG:HB2	2:C:106:HIS:NE2	2.33	0.42
1:B:291:GLU:OE1	1:B:343:HIS:CE1	2.71	0.42
3:E:11:THR:HG21	3:E:52:ASP:OD2	2.19	0.42
1:A:421:SER:O	1:A:422:GLN:HB2	2.20	0.42
2:D:176:ARG:HB2	2:D:176:ARG:NH1	2.35	0.42
1:B:82:LEU:HD23	1:B:86:LEU:HD12	2.01	0.42
2:D:17:ALA:O	2:D:21:LEU:HG	2.19	0.42
2:C:195:LEU:C	2:C:195:LEU:HD23	2.39	0.42
2:D:364:ILE:HA	2:D:368:ALA:HB3	1.99	0.42
3:F:59:LYS:HD3	3:F:59:LYS:HA	1.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:277:THR:HB	2:D:278:PRO:HD3	2.00	0.42
1:B:269:THR:HG22	3:F:148:TYR:HE1	1.84	0.42
1:A:527:ASN:O	3:E:162:ARG:NH2	2.53	0.42
1:B:442:ASN:ND2	1:B:442:ASN:C	2.73	0.42
1:A:124:LEU:HB3	1:A:137:TYR:CE1	2.55	0.42
1:B:58:GLU:HB3	1:B:132:GLU:O	2.19	0.42
2:D:42:ARG:HB2	2:D:99:ARG:HG3	2.01	0.42
1:B:403:ILE:HD12	1:B:515:LEU:CD1	2.50	0.42
2:D:365:GLU:OE1	2:D:365:GLU:CA	2.68	0.42
1:B:260:ASP:HA	1:B:261:PRO:HD3	1.94	0.42
2:C:17:ALA:O	2:C:21:LEU:HG	2.20	0.42
1:B:268:ASN:HA	1:B:268:ASN:HD22	1.65	0.42
2:D:385:LEU:O	2:D:387:GLY:N	2.53	0.42
3:E:120:PRO:HD3	3:E:128:PHE:CG	2.54	0.42
1:B:163:GLN:HG2	7:B:5009:HOH:O	2.19	0.42
2:D:169:GLU:O	2:D:170:ALA:C	2.56	0.42
1:B:90:ASN:HD22	1:B:90:ASN:HA	1.66	0.41
1:A:159:ALA:O	2:C:33:ASN:HB2	2.20	0.41
1:B:121:THR:HG21	1:B:140:GLN:CD	2.41	0.41
2:D:165:GLN:OE1	2:D:239:PHE:HA	2.20	0.41
1:B:207:VAL:HG22	1:B:313:TRP:CZ2	2.55	0.41
1:A:121:THR:HG21	1:A:140:GLN:CD	2.40	0.41
1:A:204:LEU:O	1:A:209:GLU:HG3	2.19	0.41
3:E:11:THR:HG22	3:E:12:ARG:N	2.36	0.41
1:A:330:ARG:NH1	3:E:148:TYR:HB2	2.35	0.41
2:C:263:GLU:OE2	2:C:263:GLU:HA	2.20	0.41
1:A:82:LEU:HD23	1:A:86:LEU:HD12	2.02	0.41
1:B:204:LEU:HG	1:B:205:GLN:HG3	2.02	0.41
1:A:108:ASN:HD21	1:A:175:ARG:HD3	1.86	0.41
1:B:477:GLU:C	1:B:477:GLU:OE2	2.59	0.41
2:C:270:PRO:HB3	2:D:270:PRO:CB	2.28	0.41
2:D:269:ALA:N	2:D:270:PRO:CD	2.84	0.41
1:B:123:MET:HB2	2:D:168:ARG:HD3	2.03	0.41
1:B:26:GLN:HE21	1:B:26:GLN:HB2	1.70	0.41
1:B:18:ARG:O	2:D:129:ALA:HA	2.21	0.41
1:B:249:ASN:HD22	1:B:249:ASN:HA	1.63	0.41
1:A:442:ASN:C	1:A:442:ASN:ND2	2.74	0.41
3:F:11:THR:HG22	3:F:12:ARG:N	2.37	0.40
3:E:13:ASP:HA	3:E:16:VAL:CG2	2.51	0.40
2:C:4:LEU:HD12	2:C:4:LEU:HA	1.79	0.40
1:A:246:HIS:N	1:A:246:HIS:CD2	2.89	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:22:LYS:HG3	2:C:22:LYS:O	2.21	0.40
1:B:354:TRP:CH2	1:B:499:PRO:HD3	2.56	0.40
2:C:306:ASP:O	2:C:310:SER:HB2	2.21	0.40
1:B:452:TRP:O	1:B:456:MET:HG3	2.21	0.40
2:C:364:ILE:HA	2:C:368:ALA:HB3	2.03	0.40
2:C:261:ARG:NE	2:C:285:GLN:NE2	2.63	0.40
1:A:149:HIS:CE1	2:C:105:TRP:HB2	2.56	0.40
3:E:120:PRO:HG2	3:E:125:VAL:HG12	2.04	0.40
1:B:283:THR:HB	1:B:284:PRO:HD3	2.03	0.40
1:B:114:GLU:OE1	1:B:147:HIS:CB	2.69	0.40
2:C:356:LEU:HD21	2:C:384:VAL:CG1	2.52	0.40
2:C:98:HIS:HE1	2:C:178:SER:OG	2.04	0.40
1:B:504:ASP:O	1:B:505:LYS:HB2	2.22	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	509/527 (97%)	477 (94%)	30 (6%)	2 (0%)	39	53
1	B	508/527 (96%)	479 (94%)	27 (5%)	2 (0%)	39	53
2	C	386/389 (99%)	365 (95%)	18 (5%)	3 (1%)	24	33
2	D	385/389 (99%)	360 (94%)	21 (6%)	4 (1%)	19	26
3	E	164/170 (96%)	158 (96%)	5 (3%)	1 (1%)	30	40
3	F	166/170 (98%)	158 (95%)	8 (5%)	0	100	100
All	All	2118/2172 (98%)	1997 (94%)	109 (5%)	12 (1%)	30	40

All (12) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	A	40	LYS
1	B	40	LYS
2	C	205	PRO
2	D	205	PRO
1	A	94	ARG
1	B	94	ARG
2	C	64	ALA
2	D	64	ALA
3	E	5	GLY
2	D	251	VAL
2	D	386	ALA
2	C	251	VAL

### 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	433/442 (98%)	407 (94%)	26 (6%)	24	35
1	B	432/442 (98%)	407 (94%)	25 (6%)	25	36
2	C	322/323 (100%)	310 (96%)	12 (4%)	41	60
2	D	321/323 (99%)	307 (96%)	14 (4%)	35	51
3	E	144/147 (98%)	138 (96%)	6 (4%)	36	53
3	F	146/147 (99%)	140 (96%)	6 (4%)	37	54
All	All	1798/1824 (99%)	1709 (95%)	89 (5%)	31	46

All (89) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	THR
1	A	26	GLN
1	A	30	ARG
1	A	41	ASN
1	A	43	ARG
1	A	90	ASN
1	A	110	LEU

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Mol	Chain	Res	Type
1	A	125	TRP
1	A	155	ASN
1	A	161	ASN
1	A	186	ARG
1	A	189	SER
1	A	252	GLN
1	A	269	THR
1	A	279	GLN
1	A	302	VAL
1	A	310	TYR
1	A	311	GLU
1	A	323	LYS
1	A	334	ASP
1	A	337	GLN
1	A	403	ILE
1	A	435	THR
1	A	442	ASN
1	A	477	GLU
1	A	490	SER
1	B	21	THR
1	B	26	GLN
1	B	30	ARG
1	B	41	ASN
1	B	43	ARG
1	B	90	ASN
1	B	110	LEU
1	B	125	TRP
1	B	155	ASN
1	B	161	ASN
1	B	186	ARG
1	B	252	GLN
1	B	269	THR
1	B	279	GLN
1	B	302	VAL
1	B	310	TYR
1	B	311	GLU
1	B	323	LYS
1	B	334	ASP
1	B	337	GLN
1	B	403	ILE
1	B	435	THR
1	B	442	ASN

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Mol	Chain	Res	Type
1	B	477	GLU
1	B	490	SER
2	C	80	ARG
2	C	153	LEU
2	C	168	ARG
2	C	173	ASP
2	C	179	LEU
2	C	266	GLN
2	C	311	ASP
2	C	318	ARG
2	C	356	LEU
2	C	365	GLU
2	C	378	ASP
2	C	385	LEU
2	D	35	MET
2	D	80	ARG
2	D	146	ASN
2	D	153	LEU
2	D	168	ARG
2	D	173	ASP
2	D	179	LEU
2	D	266	GLN
2	D	311	ASP
2	D	318	ARG
2	D	356	LEU
2	D	365	GLU
2	D	378	ASP
2	D	385	LEU
3	E	11	THR
3	E	16	VAL
3	E	23	ASN
3	E	44	ARG
3	E	46	SER
3	E	164	VAL
3	F	11	THR
3	F	16	VAL
3	F	23	ASN
3	F	44	ARG
3	F	46	SER
3	F	53	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (65) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	41	ASN
1	A	78	GLN
1	A	90	ASN
1	A	108	ASN
1	A	155	ASN
1	A	161	ASN
1	A	168	HIS
1	A	214	ASN
1	A	227	ASN
1	A	249	ASN
1	A	252	GLN
1	A	268	ASN
1	A	273	ASN
1	A	278	GLN
1	A	279	GLN
1	A	343	HIS
1	A	344	HIS
1	A	382	HIS
1	A	413	HIS
1	A	439	HIS
1	A	442	ASN
1	A	451	GLN
1	B	26	GLN
1	B	41	ASN
1	B	78	GLN
1	B	90	ASN
1	B	100	ASN
1	B	108	ASN
1	B	155	ASN
1	B	161	ASN
1	B	168	HIS
1	B	227	ASN
1	B	249	ASN
1	B	252	GLN
1	B	268	ASN
1	B	273	ASN
1	B	278	GLN
1	B	279	GLN
1	B	343	HIS
1	B	344	HIS
1	B	382	HIS
1	B	413	HIS

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Mol	Chain	Res	Type
1	B	439	HIS
1	B	442	ASN
1	B	451	GLN
1	B	527	ASN
2	C	98	HIS
2	C	146	ASN
2	C	161	ASN
2	C	266	GLN
2	C	285	GLN
2	D	98	HIS
2	D	146	ASN
2	D	161	ASN
2	D	266	GLN
2	D	285	GLN
3	E	39	HIS
3	E	45	ASN
3	E	144	ASN
3	E	167	GLN
3	F	39	HIS
3	F	45	ASN
3	F	144	ASN
3	F	167	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 8 ligands modelled in this entry, 7 are monoatomic - leaving 1 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	$\# Z  > 2$
6	FMT	A	6001	-	0,2,2	0.00	-	0,1,1	0.00	-

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
6	FMT	A	6001	-	-	0/0/0/0	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.

No monomer is involved in short contacts.

## 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	511/527 (96%)	0.16	25 (4%) 33 38	22, 43, 70, 86	0
1	B	510/527 (96%)	0.06	16 (3%) 52 56	23, 40, 71, 84	0
2	C	388/389 (99%)	-0.34	5 (1%) 79 81	18, 29, 47, 70	0
2	D	387/389 (99%)	0.51	35 (9%) 12 13	26, 51, 78, 86	0
3	E	166/170 (97%)	-0.18	1 (0%) 90 91	21, 34, 53, 66	0
3	F	168/170 (98%)	0.99	23 (13%) 4 5	40, 61, 82, 90	0
All	All	2130/2172 (98%)	0.15	105 (4%) 33 38	18, 41, 74, 90	0

All (105) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	205	PRO	5.8
2	D	2	SER	5.7
2	C	2	SER	5.0
2	D	353	THR	5.0
1	A	434	SER	4.5
2	D	344	ALA	4.4
2	D	375	ALA	4.3
3	F	72	PHE	4.3
2	C	205	PRO	4.1
1	A	262	ALA	4.0
1	B	244	LEU	3.9
1	B	40	LYS	3.8
1	A	326	VAL	3.8
1	A	55	GLU	3.8
2	D	343	PRO	3.8
2	C	6	GLU	3.6
3	F	27	LYS	3.5
1	A	54	ASN	3.5
1	B	55	GLU	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	261	PRO	3.5
1	B	527	ASN	3.4
1	A	310	TYR	3.4
2	D	349	LYS	3.4
1	B	333	LYS	3.4
3	F	89	SER	3.3
1	A	260	ASP	3.3
2	D	352	ILE	3.3
1	A	17	ASN	3.3
3	F	170	HIS	3.3
2	D	254	ALA	3.2
3	F	80	LYS	3.1
1	B	21	THR	3.1
3	F	26	GLU	3.1
1	B	259	ASN	3.1
1	B	258	ALA	3.0
2	D	339	PHE	2.9
3	F	102	LYS	2.9
1	B	316	ILE	2.9
2	D	260	VAL	2.8
2	D	379	GLN	2.8
1	B	59	GLN	2.8
3	E	4	LEU	2.7
2	D	256	PHE	2.7
2	D	307	PRO	2.7
1	B	19	ALA	2.7
2	D	388	LEU	2.6
3	F	67	LEU	2.6
2	D	138	PRO	2.6
2	D	44	LYS	2.6
3	F	96	ALA	2.6
3	F	83	PHE	2.5
2	D	153	LEU	2.5
2	D	45	ARG	2.5
2	C	389	LYS	2.5
2	D	357	TYR	2.5
1	A	20	PRO	2.5
3	F	21	GLN	2.5
3	F	23	ASN	2.4
1	A	320	ARG	2.4
1	A	321	LEU	2.4
2	D	356	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
2	D	193	ILE	2.4
3	F	145	LEU	2.4
2	D	252	TYR	2.4
1	B	53	ALA	2.3
3	F	22	LEU	2.3
2	D	143	GLU	2.3
2	D	220	ASN	2.3
1	B	57	LYS	2.3
2	D	374	LYS	2.3
1	B	320	ARG	2.3
1	A	366	GLN	2.3
1	A	388	GLU	2.3
3	F	30	GLU	2.3
3	F	16	VAL	2.3
1	A	263	SER	2.3
3	F	4	LEU	2.3
2	D	148	TYR	2.2
1	A	59	GLN	2.2
1	B	327	GLU	2.2
2	D	383	ALA	2.2
1	A	317	TRP	2.2
3	F	161	VAL	2.2
1	A	433	ALA	2.2
3	F	121	PRO	2.1
1	A	504	ASP	2.1
3	F	69	ALA	2.1
1	A	318	ILE	2.1
3	F	120	PRO	2.1
2	C	45	ARG	2.1
2	D	346	THR	2.1
1	A	337	GLN	2.1
2	D	308	GLU	2.1
2	D	354	ALA	2.1
1	A	259	ASN	2.1
2	D	350	GLU	2.1
2	D	26	GLU	2.1
1	A	316	ILE	2.1
2	D	259	PHE	2.1
3	F	20	ALA	2.0
1	A	338	ASP	2.0
3	F	149	ASP	2.0
1	A	324	TYR	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	525	ALA	2.0
2	D	373	PHE	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
6	FMT	A	6001	3/3	0.90	0.17	0.99	60,60,61,63	0
5	CA	A	5005	1/1	0.93	0.05	-2.42	57,57,57,57	0
4	FE	A	5002	1/1	0.97	0.04	-4.78	67,67,67,67	0
4	FE	A	5001	1/1	0.99	0.02	-6.07	47,47,47,47	0
4	FE	B	5003	1/1	0.98	0.05	-6.51	48,48,48,48	0
4	FE	B	5004	1/1	0.96	0.04	-7.09	59,59,59,59	0
5	CA	C	5007	1/1	0.84	0.06	-	72,72,72,72	0
5	CA	C	5006	1/1	0.94	0.07	-	62,62,62,62	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.