



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 02:09 am GMT

PDB ID : 4FPA  
Title : Crystal Structure of recombinant human Hexokinase type I mutant D413N  
Glucose 6-Phosphate  
Authors : Shen, L.; Honzatko, R.B.  
Deposited on : 2012-06-21  
Resolution : 2.48 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

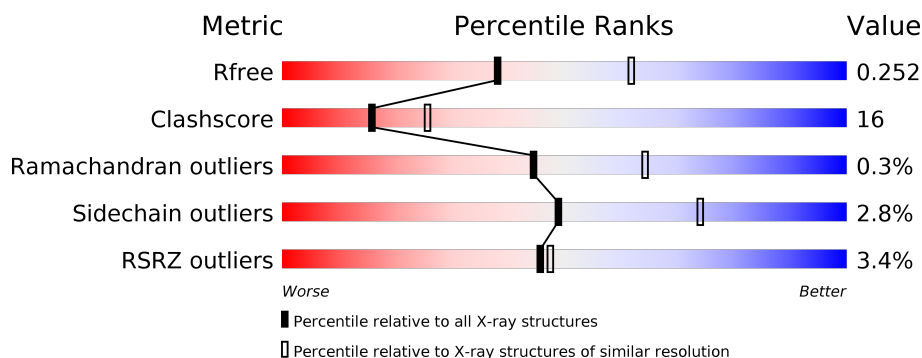
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.48 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	100719	4719 (2.50-2.46)
Clashscore	112137	5483 (2.50-2.46)
Ramachandran outliers	110173	5388 (2.50-2.46)
Sidechain outliers	110143	5390 (2.50-2.46)
RSRZ outliers	101464	4754 (2.50-2.46)

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	917	<div> <div>3%</div> <div> <div></div> <div>69%</div> <div>27%</div> <div>••</div> </div> </div>
1	B	917	<div> <div>3%</div> <div> <div></div> <div>71%</div> <div>26%</div> <div>••</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	BGC	B	1001	-	-	-	X
5	CIT	B	1007	-	-	-	X

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 14391 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.

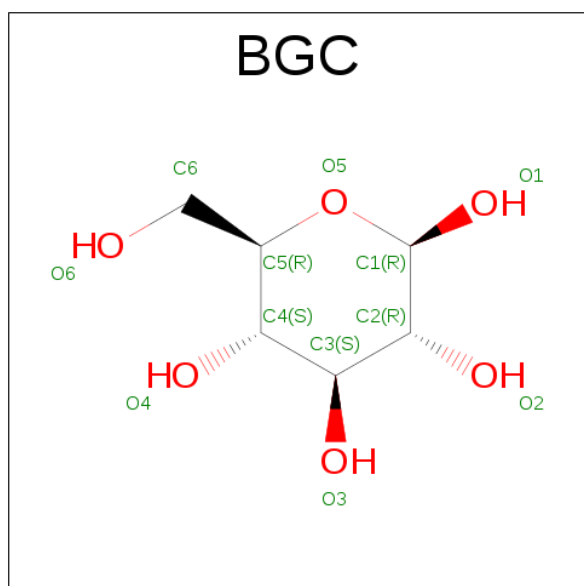
- Molecule 1 is a protein called Hexokinase-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	899	Total	C	N	O	S	0	0	0
			7032	4407	1241	1331	53			
1	B	899	Total	C	N	O	S	0	0	0
			7032	4407	1241	1331	53			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	413	ASN	ASP	ENGINEERED MUTATION	UNP P19367
B	413	ASN	ASP	ENGINEERED MUTATION	UNP P19367

- Molecule 2 is SUGAR (BETA-D-GLUCOSE) (three-letter code: BGC) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>).



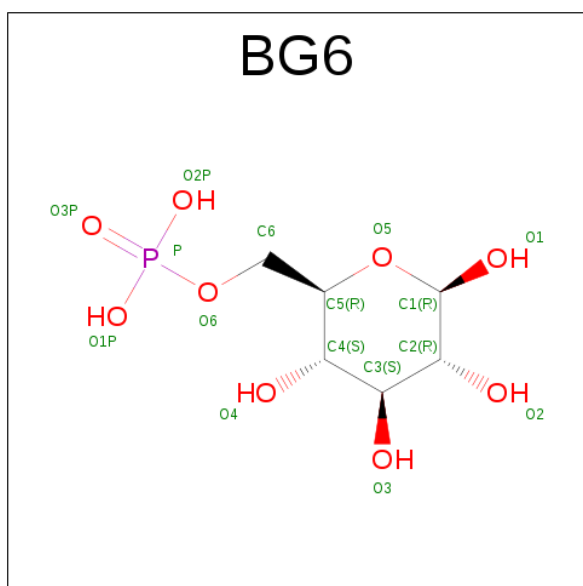
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			12	6	6		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			12	6	6		
2	B	1	Total	C	O	0	0
			12	6	6		
2	B	1	Total	C	O	0	0
			12	6	6		

- Molecule 3 is SUGAR (BETA-D-GLUCOSE-6-PHOSPHATE) (three-letter code: BG6) (formula: C<sub>6</sub>H<sub>13</sub>O<sub>9</sub>P).

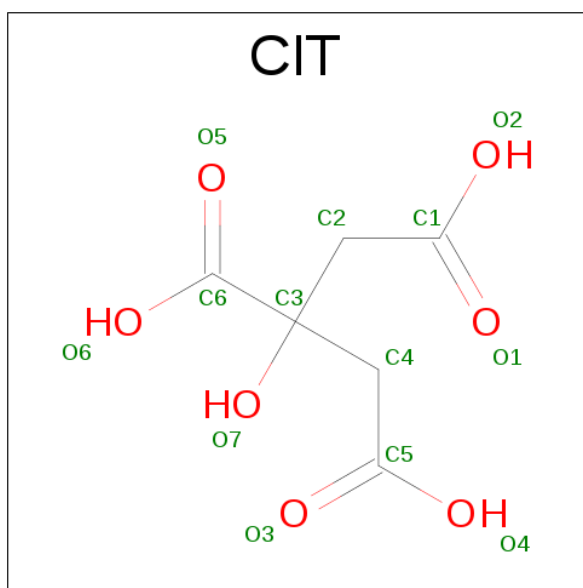


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			16	6	9	1		
3	A	1	Total	C	O	P	0	0
			16	6	9	1		
3	B	1	Total	C	O	P	0	0
			16	6	9	1		
3	B	1	Total	C	O	P	0	0
			16	6	9	1		

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

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

- Molecule 5 is CITRIC ACID (three-letter code: CIT) (formula:  $C_6H_8O_7$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			13	6	7		
5	B	1	Total	C	O	0	0
			13	6	7		

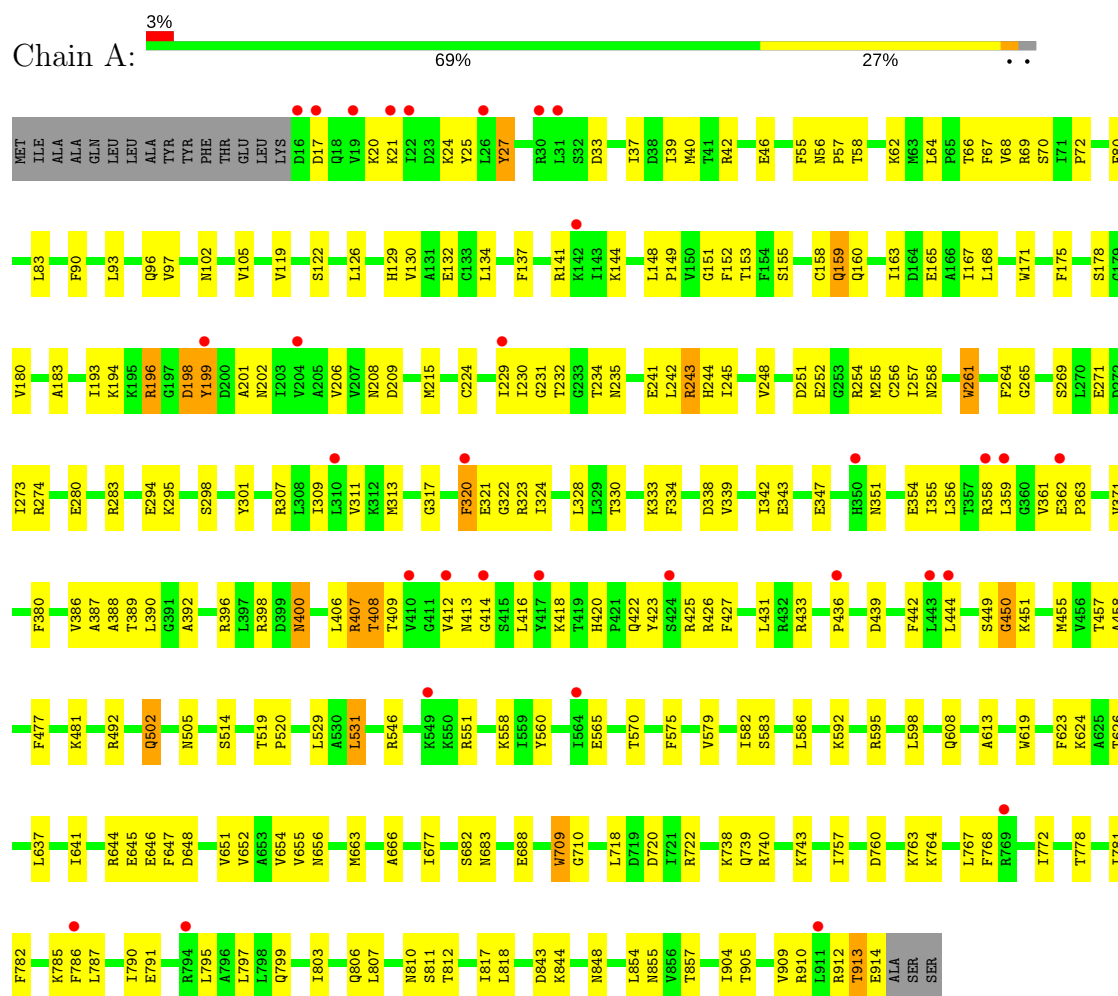
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	77	Total	O	0	0
			77	77		
6	B	108	Total	O	0	0
			108	108		

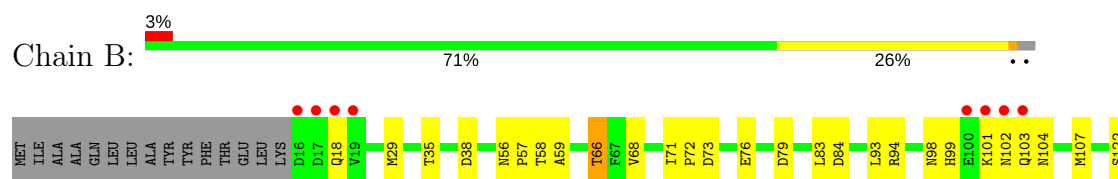
### 3 Residue-property plots

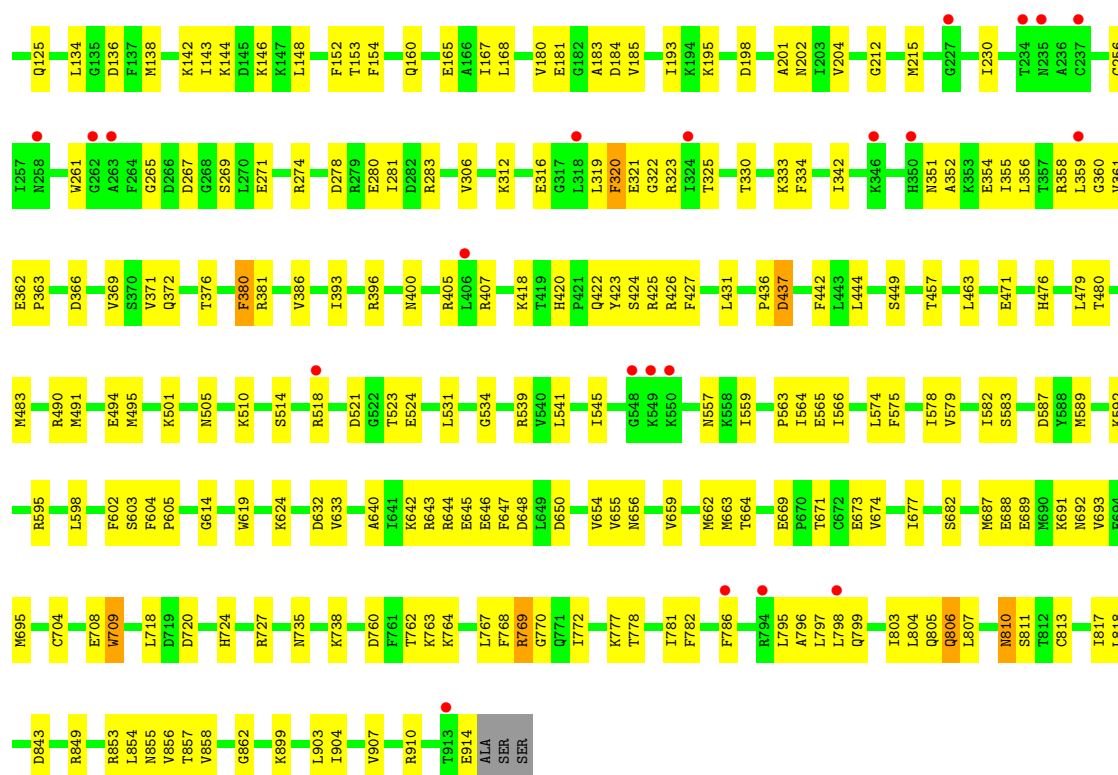
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

#### • Molecule 1: Hexokinase-1



#### • Molecule 1: Hexokinase-1







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	82.61Å 121.14Å 119.89Å 90.00° 92.69° 90.00°	Depositor
Resolution (Å)	44.15 – 2.48 44.15 – 2.48	Depositor EDS
% Data completeness (in resolution range)	98.5 (44.15-2.48) 98.5 (44.15-2.48)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.51 (at 2.48Å)	Xtriage
Refinement program	REFMAC 5.5.0066	Depositor
R, $R_{free}$	0.238 , 0.256 0.235 , 0.252	Depositor DCC
$R_{free}$ test set	4126 reflections (5.26%)	DCC
Wilson B-factor (Å <sup>2</sup> )	51.7	Xtriage
Anisotropy	0.121	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 32.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.007 for -h,-l,-k 0.000 for -h,l,k 0.019 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	14391	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.30% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA, BGC, BG6, CIT

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	0.26	0/7138	0.45	0/9606
1	B	0.25	0/7138	0.44	0/9606
All	All	0.26	0/14276	0.44	0/19212

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

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	7032	0	7092	224	0
1	B	7032	0	7092	246	0
2	A	24	0	24	2	0
2	B	24	0	24	0	0
3	A	32	0	22	2	0
3	B	32	0	22	1	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
5	A	13	0	5	1	0
5	B	13	0	5	5	0
6	A	77	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	108	0	0	6	0
All	All	14391	0	14286	465	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 16.

All (465) 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:A:132:GLU:HG3	1:A:196:ARG:NH2	1.62	1.15
1:B:797:LEU:HD11	1:B:817:ILE:HD11	1.26	1.13
1:B:518:ARG:HH12	1:B:521:ASP:HB3	1.14	1.09
1:B:518:ARG:NH1	1:B:521:ASP:HB3	1.67	1.06
1:B:480:THR:N	1:B:483:MET:HE3	1.70	1.06
1:B:66:THR:HG21	1:B:256:CYS:HB3	1.39	1.04
1:B:18:GLN:OE1	1:B:369:VAL:CG1	2.05	1.04
1:B:426:ARG:HH21	5:B:1007:CIT:H41	1.20	1.04
1:B:518:ARG:HH12	1:B:521:ASP:CB	1.73	1.02
1:B:480:THR:H	1:B:483:MET:CE	1.74	1.01
1:B:480:THR:H	1:B:483:MET:HE3	0.85	1.01
1:A:795:LEU:HD11	1:A:799:GLN:HG2	1.44	0.99
1:B:795:LEU:HD11	1:B:799:GLN:HG2	1.53	0.90
1:A:431:LEU:HD23	1:A:442:PHE:HZ	1.38	0.89
1:B:518:ARG:CZ	1:B:521:ASP:HB3	2.03	0.88
1:A:913:THR:HG23	1:A:913:THR:O	1.73	0.87
1:A:66:THR:HG22	1:A:256:CYS:O	1.76	0.86
1:B:356:LEU:HD11	1:B:371:VAL:HG21	1.55	0.85
1:B:18:GLN:OE1	1:B:369:VAL:HG12	1.76	0.85
1:B:595:ARG:NH1	1:B:650:ASP:HB2	1.93	0.83
1:A:505:ASN:HB2	6:A:1158:HOH:O	1.79	0.82
1:B:735:ASN:HB2	1:B:738:LYS:HE3	1.59	0.82
1:A:132:GLU:HG3	1:A:196:ARG:HH22	1.42	0.82
1:B:518:ARG:NH2	1:B:521:ASP:HB3	1.95	0.82
1:B:534:GLY:HA3	1:B:603:SER:HB2	1.62	0.81
1:A:361:VAL:O	1:A:363:PRO:HD3	1.83	0.79
1:A:323:ARG:NH2	1:A:362:GLU:HB2	1.98	0.79
1:B:687:MET:HE2	1:B:704:CYS:CB	2.12	0.79
1:B:539:ARG:CZ	1:B:559:ILE:HD11	2.13	0.78
1:B:564:ILE:HG23	1:B:565:GLU:N	2.00	0.77
1:B:687:MET:CE	1:B:704:CYS:HB2	2.13	0.77
1:A:767:LEU:HG	1:A:818:LEU:HD23	1.66	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:853:ARG:HB3	1:B:853:ARG:NH1	2.00	0.77
1:A:763:LYS:HG3	1:A:772:ILE:HD11	1.67	0.76
1:B:768:PHE:HA	1:B:769:ARG:HH21	1.49	0.76
1:A:786:PHE:CD2	1:A:807:LEU:HD11	2.20	0.76
1:B:575:PHE:O	1:B:579:VAL:HG23	1.86	0.76
1:A:234:THR:HG22	1:A:294:GLU:HG3	1.66	0.76
1:A:97:VAL:HG22	1:A:105:VAL:HA	1.67	0.75
1:B:18:GLN:NE2	1:B:366:ASP:O	2.18	0.75
1:B:778:THR:O	1:B:781:ILE:HG12	1.87	0.75
1:B:18:GLN:OE1	1:B:369:VAL:CB	2.35	0.74
1:B:66:THR:CG2	1:B:256:CYS:HB3	2.17	0.74
1:B:642:LYS:O	1:B:645:GLU:HG2	1.88	0.74
1:B:18:GLN:OE1	1:B:369:VAL:HB	1.87	0.74
1:A:786:PHE:CZ	1:A:790:ILE:HG13	2.23	0.73
1:B:767:LEU:HG	1:B:818:LEU:HD23	1.70	0.73
1:B:518:ARG:HH22	1:B:521:ASP:HB3	1.52	0.73
1:B:797:LEU:HD11	1:B:817:ILE:CD1	2.13	0.73
1:B:29:MET:HE1	1:B:381:ARG:CZ	2.18	0.73
1:B:35:THR:O	1:B:38:ASP:HB3	1.88	0.73
1:A:160:GLN:HG2	1:A:165:GLU:O	1.89	0.72
1:B:518:ARG:HH22	1:B:521:ASP:CB	2.01	0.72
1:B:431:LEU:HD22	1:B:442:PHE:HZ	1.53	0.72
1:A:502:GLN:OE1	1:A:502:GLN:N	2.17	0.71
1:B:687:MET:HE1	1:B:704:CYS:HB2	1.71	0.71
1:A:209:ASP:CG	1:A:229:ILE:HD12	2.11	0.71
1:B:853:ARG:HH11	1:B:853:ARG:HB3	1.54	0.71
1:B:688:GLU:OE1	1:B:689:GLU:OE1	2.07	0.71
1:B:420:HIS:HD2	1:B:423:TYR:H	1.39	0.70
1:A:570:THR:HA	1:A:626:THR:OG1	1.92	0.70
1:B:323:ARG:CZ	1:B:362:GLU:HB2	2.22	0.70
1:B:323:ARG:NH1	1:B:362:GLU:HB2	2.07	0.70
1:A:144:LYS:NZ	1:A:198:ASP:OD1	2.25	0.69
1:A:39:ILE:HD13	1:A:42:ARG:HH21	1.57	0.69
1:A:57:PRO:HG2	1:B:799:GLN:HE21	1.58	0.69
1:A:354:GLU:O	1:A:358:ARG:HG3	1.92	0.69
1:A:323:ARG:CZ	1:A:362:GLU:HB2	2.23	0.68
1:A:159:GLN:HB3	1:A:167:ILE:HB	1.75	0.68
1:A:431:LEU:HD23	1:A:442:PHE:CZ	2.25	0.68
1:A:431:LEU:CD2	1:A:442:PHE:HZ	2.07	0.68
1:A:398:ARG:HB3	1:A:398:ARG:HH11	1.59	0.68
1:A:799:GLN:NE2	1:B:58:THR:HB	2.08	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:267:ASP:OD1	1:B:269:SER:HB2	1.94	0.68
1:A:152:PHE:HB3	1:A:206:VAL:HG22	1.74	0.68
1:B:687:MET:CE	1:B:704:CYS:CB	2.72	0.67
1:A:56:ASN:N	1:A:57:PRO:HD2	2.09	0.67
1:B:122:SER:OG	1:B:125:GLN:HG3	1.94	0.67
1:B:160:GLN:HG2	1:B:165:GLU:O	1.93	0.67
1:B:351:ASN:O	1:B:355:ILE:HG12	1.95	0.67
1:B:810:ASN:HD22	1:B:810:ASN:H	1.43	0.66
1:A:241:GLU:O	1:A:245:ILE:HD12	1.95	0.66
1:B:772:ILE:HG22	1:B:777:LYS:HD2	1.78	0.66
1:A:132:GLU:CG	1:A:196:ARG:NH2	2.52	0.66
1:A:230:ILE:HD11	1:A:386:VAL:HG11	1.77	0.66
1:A:317:GLY:HA2	1:A:321:GLU:O	1.94	0.66
1:A:778:THR:HB	1:A:781:ILE:HD13	1.75	0.66
1:B:98:ASN:HB3	1:B:103:GLN:HB2	1.78	0.66
1:B:142:LYS:HE2	1:B:142:LYS:HA	1.78	0.66
1:B:564:ILE:HG23	1:B:565:GLU:H	1.60	0.66
1:B:760:ASP:O	1:B:764:LYS:HG2	1.96	0.66
1:A:245:ILE:HG12	1:A:257:ILE:HD11	1.77	0.66
1:B:425:ARG:HH22	5:B:1007:CIT:H42	1.61	0.66
1:A:280:GLU:HG3	1:A:283:ARG:NH2	2.11	0.65
1:A:309:ILE:O	1:A:313:MET:HG3	1.96	0.65
1:A:913:THR:O	1:A:913:THR:CG2	2.45	0.65
1:A:619:TRP:HB3	1:A:623:PHE:O	1.95	0.65
1:B:644:ARG:NE	1:B:646:GLU:OE1	2.29	0.65
1:A:40:MET:HG3	1:A:388:ALA:O	1.96	0.65
1:A:407:ARG:HG2	1:A:439:ASP:HB2	1.78	0.65
1:B:431:LEU:CD2	1:B:442:PHE:HZ	2.10	0.65
1:A:93:LEU:HD22	1:A:450:GLY:HA3	1.79	0.65
1:B:405:ARG:HD2	1:B:437:ASP:HB3	1.79	0.64
1:A:24:LYS:O	1:A:27:TYR:HB3	1.97	0.64
1:A:356:LEU:HD11	1:A:371:VAL:HG21	1.80	0.64
1:B:693:VAL:HG12	1:B:693:VAL:O	1.98	0.64
1:B:405:ARG:HB3	1:B:405:ARG:HH11	1.63	0.64
1:B:405:ARG:HB3	1:B:405:ARG:NH1	2.13	0.64
1:B:539:ARG:HB3	1:B:559:ILE:HD13	1.79	0.63
1:A:196:ARG:HG2	1:A:198:ASP:HB2	1.80	0.63
1:A:663:MET:HG3	1:A:904:ILE:HD11	1.81	0.63
1:B:423:TYR:HB3	6:B:1181:HOH:O	1.99	0.63
1:B:134:LEU:O	1:B:138:MET:HG3	1.99	0.62
1:B:810:ASN:HD22	1:B:810:ASN:N	1.97	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:583:SER:OG	1:B:592:LYS:NZ	2.32	0.62
1:A:320:PHE:O	1:A:321:GLU:HB2	1.99	0.62
1:B:813:CYS:O	1:B:817:ILE:HG12	1.98	0.62
1:A:420:HIS:HD2	1:A:423:TYR:N	1.98	0.61
1:B:426:ARG:HE	5:B:1007:CIT:H22	1.65	0.61
1:A:66:THR:CG2	1:A:256:CYS:HB3	2.30	0.61
1:A:799:GLN:HE22	1:B:58:THR:HB	1.65	0.61
1:B:854:LEU:HD12	1:B:855:ASN:N	2.15	0.61
1:A:795:LEU:HD11	1:A:799:GLN:CG	2.26	0.61
1:B:101:LYS:C	1:B:103:GLN:H	2.04	0.61
1:A:767:LEU:HD13	1:A:768:PHE:CE2	2.35	0.61
1:B:18:GLN:OE1	1:B:369:VAL:HG11	1.99	0.60
1:B:29:MET:HE1	1:B:381:ARG:NH2	2.15	0.60
1:B:687:MET:HE2	1:B:704:CYS:HA	1.83	0.60
1:B:768:PHE:CE1	1:B:811:SER:HB3	2.37	0.60
1:A:558:LYS:HD3	1:A:560:TYR:OH	2.01	0.60
1:B:782:PHE:HD1	1:B:786:PHE:CE1	2.19	0.60
1:B:687:MET:HE2	1:B:704:CYS:CA	2.31	0.60
1:B:98:ASN:ND2	1:B:101:LYS:HD2	2.17	0.59
1:A:912:ARG:C	1:A:914:GLU:H	2.06	0.59
1:B:426:ARG:HH21	5:B:1007:CIT:C4	2.06	0.59
1:B:763:LYS:HG3	1:B:772:ILE:HD11	1.83	0.59
1:B:768:PHE:HE1	1:B:811:SER:HB3	1.65	0.59
1:B:66:THR:HG23	1:B:68:VAL:H	1.68	0.59
1:A:910:ARG:O	1:A:914:GLU:HB2	2.02	0.59
1:B:471:GLU:HG3	6:B:1161:HOH:O	2.02	0.59
1:A:398:ARG:HB3	1:A:398:ARG:NH1	2.16	0.59
1:B:564:ILE:CG2	1:B:565:GLU:N	2.65	0.59
1:B:18:GLN:HE22	1:B:366:ASP:C	2.06	0.59
1:A:209:ASP:OD1	1:A:229:ILE:HD12	2.03	0.59
1:B:58:THR:HG22	1:B:58:THR:O	2.03	0.58
1:A:193:ILE:HD13	1:A:201:ALA:HB3	1.86	0.58
1:B:193:ILE:HD13	1:B:201:ALA:HB3	1.86	0.58
1:A:644:ARG:NE	1:A:646:GLU:OE1	2.28	0.58
1:A:361:VAL:C	1:A:363:PRO:HD3	2.24	0.57
1:B:564:ILE:CG2	1:B:565:GLU:H	2.17	0.57
1:A:546:ARG:O	1:A:551:ARG:HA	2.05	0.57
1:B:479:LEU:HA	1:B:483:MET:HE1	1.85	0.57
1:B:619:TRP:CD1	1:B:624:LYS:HA	2.40	0.57
1:B:56:ASN:N	1:B:57:PRO:HD2	2.19	0.57
1:B:605:PRO:HB3	1:B:708:GLU:HG3	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:595:ARG:HD2	1:A:648:ASP:OD2	2.05	0.56
1:B:426:ARG:NH2	5:B:1007:CIT:H41	2.04	0.56
1:A:389:THR:O	1:A:392:ALA:HB3	2.05	0.56
1:B:265:GLY:HA2	1:B:269:SER:HB3	1.87	0.56
1:A:235:ASN:HA	1:A:261:TRP:CD1	2.40	0.56
1:A:768:PHE:HE1	1:A:811:SER:HB3	1.71	0.56
1:B:644:ARG:HD3	1:B:646:GLU:OE1	2.05	0.56
1:B:687:MET:HE2	1:B:704:CYS:HB2	1.82	0.56
1:A:271:GLU:OE1	1:A:271:GLU:HA	2.04	0.56
1:A:66:THR:HG21	1:A:256:CYS:HB3	1.87	0.55
1:A:307:ARG:O	1:A:311:VAL:HG23	2.05	0.55
1:B:420:HIS:HD2	1:B:423:TYR:N	2.02	0.55
1:A:155:SER:HA	1:A:208:ASN:ND2	2.20	0.55
1:A:760:ASP:O	1:A:764:LYS:HG2	2.05	0.55
1:A:688:GLU:OE1	1:A:848:ASN:ND2	2.38	0.55
1:B:230:ILE:HD11	1:B:386:VAL:HG11	1.87	0.55
1:B:724:HIS:HA	1:B:727:ARG:NH1	2.21	0.55
1:B:565:GLU:CG	1:B:566:ILE:N	2.70	0.55
1:A:171:TRP:HB3	1:A:175:PHE:O	2.06	0.55
1:B:342:ILE:HG22	1:B:342:ILE:O	2.06	0.55
1:A:126:LEU:O	1:A:129:HIS:HB3	2.07	0.55
1:B:76:GLU:H	1:B:76:GLU:CD	2.10	0.55
1:B:565:GLU:HG2	1:B:566:ILE:H	1.70	0.55
1:A:558:LYS:HD3	1:A:560:TYR:CZ	2.42	0.54
1:B:101:LYS:O	1:B:103:GLN:N	2.40	0.54
1:A:420:HIS:HD2	1:A:423:TYR:H	1.53	0.54
1:A:320:PHE:C	1:A:322:GLY:H	2.10	0.54
1:B:59:ALA:HB1	6:B:1172:HOH:O	2.07	0.54
1:A:739:GLN:O	1:A:743:LYS:HG3	2.07	0.54
1:A:418:LYS:HG3	1:A:444:LEU:HD11	1.88	0.54
1:A:72:PRO:HD3	1:A:215:MET:CE	2.38	0.54
1:B:786:PHE:CE2	1:B:803:ILE:HG22	2.42	0.54
1:A:168:LEU:HD23	1:A:180:VAL:HG12	1.89	0.54
1:A:380:PHE:HD2	1:A:426:ARG:HD3	1.72	0.54
1:B:602:PHE:CE2	1:B:633:VAL:HG11	2.42	0.54
1:A:244:HIS:HD2	1:A:396:ARG:NH2	2.06	0.53
1:A:608:GLN:OE1	1:A:654:VAL:HG13	2.08	0.53
1:B:418:LYS:HG2	1:B:444:LEU:HD21	1.90	0.53
1:B:479:LEU:HA	1:B:483:MET:CE	2.38	0.53
1:A:422:GLN:OE1	1:A:425:ARG:NH2	2.37	0.53
1:A:575:PHE:O	1:A:579:VAL:HG23	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:212:GLY:HA3	1:B:449:SER:HB2	1.89	0.53
1:A:58:THR:OG1	1:B:799:GLN:NE2	2.42	0.53
1:B:565:GLU:CG	1:B:566:ILE:H	2.22	0.53
1:B:910:ARG:HG2	1:B:914:GLU:OE2	2.09	0.53
1:A:502:GLN:CD	1:A:502:GLN:H	2.09	0.53
1:A:134:LEU:HD23	1:A:199:TYR:HE1	1.71	0.53
1:A:738:LYS:HE2	6:A:1148:HOH:O	2.08	0.53
1:B:505:ASN:HB2	6:B:1185:HOH:O	2.09	0.53
1:B:29:MET:CE	1:B:381:ARG:CZ	2.86	0.53
1:A:351:ASN:O	1:A:355:ILE:HG12	2.10	0.52
1:B:735:ASN:HB2	1:B:738:LYS:CE	2.33	0.52
1:B:673:GLU:CD	1:B:849:ARG:HH22	2.13	0.52
1:A:155:SER:HA	1:A:208:ASN:HD22	1.74	0.52
1:A:122:SER:HA	1:A:178:SER:OG	2.10	0.52
1:B:393:ILE:O	1:B:396:ARG:HB3	2.08	0.52
1:A:324:ILE:HG23	1:A:328:LEU:HD23	1.90	0.52
1:B:319:LEU:HB3	1:B:320:PHE:CE1	2.45	0.52
1:B:491:MET:O	1:B:495:MET:HG3	2.09	0.52
1:A:451:LYS:O	1:A:455:MET:HG2	2.09	0.52
1:A:420:HIS:CD2	1:A:423:TYR:HB2	2.44	0.52
1:B:354:GLU:O	1:B:358:ARG:HG3	2.10	0.52
1:B:687:MET:CE	1:B:704:CYS:SG	2.98	0.52
1:A:646:GLU:HB2	1:A:647:PHE:CD2	2.45	0.51
1:A:317:GLY:CA	1:A:321:GLU:O	2.57	0.51
1:B:514:SER:OG	1:B:704:CYS:HB3	2.10	0.51
1:A:387:ALA:HA	1:A:427:PHE:HE1	1.76	0.51
1:B:644:ARG:CD	1:B:646:GLU:OE1	2.58	0.51
1:A:137:PHE:O	1:A:141:ARG:HD3	2.10	0.51
1:B:574:LEU:O	1:B:578:ILE:HG12	2.11	0.51
1:A:418:LYS:HG2	1:A:444:LEU:HD21	1.92	0.51
1:B:674:VAL:HB	1:B:858:VAL:HG22	1.92	0.51
1:B:531:LEU:HD21	1:B:582:ILE:HD11	1.91	0.51
1:B:763:LYS:CG	1:B:772:ILE:HD11	2.41	0.51
1:A:307:ARG:HB2	1:A:334:PHE:HB3	1.93	0.51
1:A:224:CYS:HA	1:A:409:THR:O	2.11	0.51
1:A:763:LYS:CG	1:A:772:ILE:HD11	2.39	0.51
1:B:306:VAL:HG11	1:B:334:PHE:CE2	2.46	0.50
1:A:235:ASN:HA	1:A:261:TRP:NE1	2.25	0.50
1:A:33:ASP:O	1:A:37:ILE:HG12	2.12	0.50
1:A:69:ARG:O	1:A:70:SER:HB3	2.12	0.50
1:B:782:PHE:CD1	1:B:786:PHE:CE1	2.99	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93:LEU:HD12	1:B:93:LEU:N	2.26	0.50
1:A:655:VAL:HG12	1:A:656:ASN:O	2.12	0.50
1:B:278:ASP:O	1:B:281:ILE:HG22	2.11	0.50
1:A:414:GLY:HA2	3:A:1002:BG6:O6	2.11	0.50
1:A:787:LEU:O	1:A:791:GLU:HG3	2.11	0.50
1:A:646:GLU:HB2	1:A:647:PHE:HD2	1.76	0.50
1:A:93:LEU:N	1:A:93:LEU:HD12	2.27	0.50
1:A:97:VAL:HG21	1:A:105:VAL:HG22	1.93	0.50
1:B:687:MET:HE2	1:B:704:CYS:SG	2.52	0.50
1:B:786:PHE:HE2	1:B:803:ILE:HG22	1.76	0.50
1:A:265:GLY:HA2	1:A:269:SER:OG	2.12	0.50
1:A:514:SER:HA	1:A:608:GLN:HE22	1.76	0.50
1:A:67:PHE:HA	1:A:255:MET:SD	2.52	0.50
1:B:646:GLU:HB3	1:B:647:PHE:HD2	1.77	0.50
1:B:563:PRO:HG2	1:B:566:ILE:HG13	1.94	0.49
1:A:97:VAL:CG2	1:A:105:VAL:HG22	2.41	0.49
1:B:662:MET:SD	1:B:687:MET:HE3	2.53	0.49
1:B:98:ASN:O	1:B:103:GLN:O	2.30	0.49
1:A:245:ILE:HG22	1:A:245:ILE:O	2.12	0.49
1:A:786:PHE:CE2	1:A:807:LEU:HD11	2.48	0.49
1:B:168:LEU:HD23	1:B:180:VAL:HG12	1.95	0.49
1:B:524:GLU:H	1:B:524:GLU:CD	2.16	0.49
1:B:79:ASP:HB3	1:B:148:LEU:CD2	2.43	0.49
1:A:248:VAL:HG21	1:A:255:MET:HE1	1.94	0.49
1:A:598:LEU:HD23	1:A:598:LEU:C	2.33	0.48
1:A:356:LEU:HD23	1:A:359:LEU:HD12	1.94	0.48
1:A:412:VAL:HG12	1:A:413:ASN:N	2.27	0.48
1:A:163:ILE:HG13	1:A:163:ILE:O	2.13	0.48
1:A:141:ARG:HD2	1:A:141:ARG:N	2.28	0.48
1:B:79:ASP:HB3	1:B:148:LEU:HD22	1.96	0.48
1:A:242:LEU:C	1:A:244:HIS:H	2.16	0.48
1:B:677:ILE:O	1:B:682:SER:HA	2.14	0.48
1:A:637:LEU:O	1:A:641:ILE:HG13	2.14	0.48
1:A:782:PHE:CD1	1:A:786:PHE:CE1	3.02	0.48
1:B:136:ASP:C	1:B:136:ASP:OD1	2.51	0.48
1:A:198:ASP:HB3	1:A:199:TYR:HD2	1.79	0.48
1:B:518:ARG:HH12	1:B:521:ASP:CA	2.25	0.48
1:B:523:THR:OG1	1:B:910:ARG:NH1	2.47	0.48
1:B:669:GLU:OE2	1:B:671:THR:N	2.40	0.48
1:B:342:ILE:O	1:B:372:GLN:HG3	2.13	0.47
1:A:320:PHE:O	1:A:322:GLY:N	2.40	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:608:GLN:HG2	1:A:613:ALA:O	2.15	0.47
1:B:66:THR:CG2	1:B:68:VAL:H	2.25	0.47
1:B:654:VAL:HG13	1:B:654:VAL:O	2.13	0.47
1:A:905:THR:O	1:A:909:VAL:HG23	2.14	0.47
1:A:321:GLU:N	1:A:321:GLU:OE1	2.47	0.47
1:B:271:GLU:OE2	1:B:274:ARG:HD3	2.14	0.47
1:B:58:THR:CG2	1:B:58:THR:O	2.63	0.47
1:A:83:LEU:O	1:A:153:THR:HB	2.15	0.47
1:A:66:THR:O	1:A:67:PHE:HB2	2.15	0.47
1:B:671:THR:OG1	1:B:857:THR:HG23	2.15	0.47
1:B:786:PHE:CE2	1:B:807:LEU:HD11	2.49	0.47
1:A:66:THR:HG23	1:A:68:VAL:H	1.79	0.47
1:B:598:LEU:HD23	1:B:598:LEU:C	2.35	0.47
1:B:796:ALA:O	1:B:799:GLN:HB3	2.14	0.47
1:A:519:THR:HB	1:A:520:PRO:HD2	1.96	0.47
1:B:420:HIS:CD2	1:B:423:TYR:H	2.26	0.47
1:B:640:ALA:N	1:B:643:ARG:HH21	2.13	0.47
1:A:416:LEU:HD21	1:A:423:TYR:CE2	2.49	0.46
1:A:390:LEU:HD23	1:A:431:LEU:HD22	1.98	0.46
1:A:654:VAL:O	1:A:654:VAL:HG13	2.15	0.46
1:B:664:THR:HG23	1:B:899:LYS:HB3	1.98	0.46
1:B:689:GLU:N	1:B:689:GLU:OE1	2.48	0.46
1:A:652:VAL:HG21	1:A:909:VAL:HG22	1.97	0.46
1:B:797:LEU:HD21	1:B:817:ILE:HD13	1.97	0.46
1:A:529:LEU:HD11	1:A:586:LEU:HD21	1.97	0.46
1:B:810:ASN:N	1:B:810:ASN:ND2	2.63	0.46
1:A:56:ASN:N	1:A:57:PRO:CD	2.77	0.46
1:A:583:SER:OG	1:A:592:LYS:NZ	2.49	0.46
1:A:785:LYS:HB2	6:A:1159:HOH:O	2.15	0.46
1:B:101:LYS:C	1:B:103:GLN:N	2.68	0.46
1:B:342:ILE:CG2	1:B:342:ILE:O	2.63	0.46
1:A:408:THR:OG1	1:A:409:THR:N	2.49	0.46
1:B:769:ARG:HH22	1:B:811:SER:HA	1.81	0.46
1:A:90:PHE:CZ	1:A:130:VAL:HG13	2.51	0.46
1:A:24:LYS:O	1:A:27:TYR:CB	2.63	0.46
1:A:64:LEU:HD13	1:A:158:CYS:O	2.16	0.46
1:B:165:GLU:HG3	1:B:184:ASP:OD2	2.16	0.46
1:A:295:LYS:HA	1:A:301:TYR:CD2	2.51	0.46
1:B:856:VAL:HG22	1:B:857:THR:N	2.31	0.46
1:A:339:VAL:O	1:A:343:GLU:HG3	2.16	0.45
1:B:376:THR:O	1:B:380:PHE:HB2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:691:LYS:HG3	1:B:692:ASN:N	2.31	0.45
1:A:425:ARG:NH1	5:A:1007:CIT:O2	2.50	0.45
1:A:803:ILE:O	1:A:806:GLN:HB2	2.16	0.45
1:B:320:PHE:N	1:B:320:PHE:CD1	2.85	0.45
1:A:398:ARG:CB	1:A:398:ARG:HH11	2.27	0.45
1:A:799:GLN:NE2	1:B:58:THR:CB	2.78	0.45
1:A:193:ILE:O	1:A:194:LYS:C	2.55	0.45
1:B:181:GLU:HG3	6:B:1189:HOH:O	2.16	0.45
1:B:195:LYS:HG2	1:B:195:LYS:O	2.17	0.45
1:B:539:ARG:HG3	1:B:541:LEU:HD11	1.99	0.45
1:B:589:MET:CE	1:B:589:MET:HA	2.46	0.45
1:B:595:ARG:NE	1:B:648:ASP:OD2	2.50	0.45
1:B:320:PHE:C	1:B:322:GLY:H	2.19	0.45
1:B:659:VAL:HG22	1:B:704:CYS:SG	2.56	0.45
1:B:718:LEU:C	1:B:720:ASP:N	2.70	0.45
1:A:320:PHE:C	1:A:322:GLY:N	2.71	0.45
1:B:361:VAL:HG23	1:B:363:PRO:HD3	1.98	0.45
1:B:646:GLU:HB3	1:B:647:PHE:CD2	2.51	0.45
1:A:514:SER:HA	1:A:608:GLN:NE2	2.32	0.45
1:A:62:LYS:HB3	1:A:64:LEU:CD2	2.46	0.45
1:A:232:THR:O	1:A:298:SER:HB2	2.17	0.44
1:B:862:GLY:HA2	3:B:1004:BG6:O6	2.17	0.44
1:A:193:ILE:O	1:A:196:ARG:N	2.49	0.44
1:A:134:LEU:CD2	1:A:199:TYR:HE1	2.31	0.44
1:B:539:ARG:HB3	1:B:559:ILE:CD1	2.44	0.44
1:A:722:ARG:CZ	1:A:740:ARG:HD3	2.46	0.44
1:B:330:THR:HB	1:B:333:LYS:HG3	1.98	0.44
1:B:480:THR:HG23	1:B:483:MET:CE	2.47	0.44
1:A:168:LEU:CD2	1:A:180:VAL:HG12	2.48	0.44
1:B:204:VAL:HG12	1:B:457:THR:HG23	2.00	0.44
1:A:323:ARG:NH2	1:A:362:GLU:O	2.51	0.44
1:A:46:GLU:HG3	1:A:264:PHE:CE1	2.53	0.44
1:A:786:PHE:CE1	1:A:790:ILE:HG13	2.53	0.44
1:B:144:LYS:NZ	1:B:198:ASP:HB3	2.33	0.44
1:A:252:GLU:OE2	1:A:812:THR:HB	2.17	0.44
1:B:71:ILE:HB	1:B:72:PRO:HD2	2.00	0.44
1:B:762:THR:O	1:B:770:GLY:HA2	2.17	0.44
1:A:119:VAL:O	1:A:119:VAL:HG12	2.18	0.44
1:A:39:ILE:HD11	1:A:273:ILE:HG12	2.00	0.44
1:A:80:PHE:CE2	1:A:458:ALA:HA	2.53	0.44
1:B:280:GLU:HG3	1:B:283:ARG:NH2	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:718:LEU:C	1:A:720:ASP:N	2.70	0.43
1:A:251:ASP:N	1:A:251:ASP:OD1	2.51	0.43
1:A:323:ARG:HH21	1:A:362:GLU:HB2	1.80	0.43
1:A:797:LEU:HD21	1:A:817:ILE:HG12	2.00	0.43
1:B:614:GLY:O	1:B:632:ASP:HA	2.18	0.43
1:A:595:ARG:CD	1:A:648:ASP:OD2	2.67	0.43
1:B:420:HIS:CD2	1:B:423:TYR:HB2	2.53	0.43
1:B:579:VAL:HG21	1:B:640:ALA:CB	2.48	0.43
1:A:244:HIS:CD2	1:A:396:ARG:NH2	2.85	0.43
1:A:406:LEU:HD12	1:A:407:ARG:H	1.84	0.43
1:A:477:PHE:CZ	1:A:757:ILE:HD11	2.53	0.43
1:A:677:ILE:HG21	2:A:1003:BGC:C6	2.48	0.43
1:B:146:LYS:HB2	1:B:148:LEU:HG	2.00	0.43
1:B:101:LYS:O	1:B:103:GLN:CG	2.66	0.43
1:A:148:LEU:HA	1:A:149:PRO:HD3	1.82	0.43
1:A:413:ASN:ND2	3:A:1002:BG6:O1	2.45	0.43
1:B:386:VAL:HG12	1:B:427:PHE:CE1	2.54	0.43
1:A:519:THR:HB	1:A:666:ALA:HB1	2.00	0.43
1:B:154:PHE:CE2	1:B:185:VAL:HG11	2.54	0.43
1:A:400:ASN:HD22	1:A:400:ASN:HA	1.60	0.43
1:A:677:ILE:O	1:A:682:SER:HA	2.19	0.43
1:B:644:ARG:HD3	1:B:646:GLU:HB2	2.01	0.43
1:B:655:VAL:HG12	1:B:656:ASN:O	2.19	0.43
1:B:907:VAL:HG12	1:B:907:VAL:O	2.19	0.43
1:A:25:TYR:C	1:A:27:TYR:H	2.21	0.42
1:B:101:LYS:O	1:B:103:GLN:HG3	2.19	0.42
1:B:640:ALA:HA	1:B:643:ARG:HE	1.84	0.42
1:A:274:ARG:NH1	6:A:1156:HOH:O	2.52	0.42
1:A:449:SER:O	1:A:450:GLY:C	2.58	0.42
1:A:912:ARG:C	1:A:914:GLU:N	2.72	0.42
1:B:84:ASP:HA	1:B:153:THR:HB	2.01	0.42
1:B:323:ARG:NH2	1:B:362:GLU:HB2	2.33	0.42
1:B:640:ALA:CA	1:B:643:ARG:HH21	2.31	0.42
1:A:199:TYR:CD2	1:A:199:TYR:N	2.87	0.42
1:A:330:THR:HB	1:A:333:LYS:HG3	2.01	0.42
1:B:541:LEU:HG	1:B:557:ASN:CB	2.48	0.42
1:A:55:PHE:C	1:A:57:PRO:HD2	2.39	0.42
1:A:677:ILE:HG21	2:A:1003:BGC:H6C1	2.00	0.42
1:A:492:ARG:CZ	1:A:844:LYS:HG3	2.49	0.42
1:B:640:ALA:HB2	1:B:643:ARG:NH2	2.35	0.42
1:A:433:ARG:O	1:A:436:PRO:HD3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:644:ARG:O	1:A:645:GLU:HB3	2.20	0.42
1:B:420:HIS:CD2	1:B:422:GLN:H	2.38	0.42
1:A:230:ILE:HD13	1:A:386:VAL:HG21	2.00	0.42
1:A:258:ASN:OD1	1:A:258:ASN:C	2.57	0.42
1:B:66:THR:HG23	1:B:68:VAL:HG23	2.02	0.42
1:B:167:ILE:HA	1:B:183:ALA:O	2.20	0.42
1:A:231:GLY:O	1:A:232:THR:C	2.58	0.42
1:A:420:HIS:CD2	1:A:422:GLN:H	2.37	0.42
1:A:619:TRP:CD1	1:A:624:LYS:HA	2.55	0.42
1:A:595:ARG:NE	1:A:648:ASP:OD2	2.52	0.42
1:B:587:ASP:OD1	1:B:592:LYS:HE2	2.19	0.42
1:B:579:VAL:HG21	1:B:640:ALA:HB3	2.01	0.42
1:B:501:LYS:HB3	1:B:695:MET:SD	2.60	0.42
1:A:134:LEU:HD23	1:A:199:TYR:CE1	2.52	0.42
1:B:640:ALA:HA	1:B:643:ARG:NE	2.34	0.42
1:B:854:LEU:HD12	1:B:855:ASN:H	1.82	0.42
1:B:356:LEU:HD23	1:B:359:LEU:HD12	2.02	0.41
1:B:325:THR:HG21	1:B:360:GLY:HA3	2.00	0.41
1:B:510:LYS:HA	1:B:510:LYS:HD3	1.92	0.41
1:B:604:PHE:HB3	1:B:605:PRO:HD2	2.01	0.41
1:B:94:ARG:NH1	1:B:143:ILE:HG21	2.35	0.41
1:A:105:VAL:HG11	1:A:451:LYS:HE2	2.02	0.41
1:A:20:LYS:O	1:A:24:LYS:HG3	2.20	0.41
1:A:492:ARG:O	1:A:492:ARG:HD2	2.21	0.41
1:A:347:GLU:HB3	1:A:351:ASN:ND2	2.35	0.41
1:A:151:GLY:HA3	1:A:457:THR:OG1	2.19	0.41
1:A:710:GLY:O	1:A:739:GLN:HA	2.21	0.41
1:B:72:PRO:HD3	1:B:215:MET:CE	2.50	0.41
1:B:73:ASP:OD1	1:B:73:ASP:C	2.59	0.41
1:A:854:LEU:HD12	1:A:855:ASN:N	2.36	0.41
1:B:431:LEU:HD22	1:B:442:PHE:CZ	2.43	0.41
1:A:167:ILE:HA	1:A:183:ALA:O	2.21	0.41
1:B:476:HIS:CD2	6:B:1194:HOH:O	2.74	0.41
1:A:72:PRO:HG3	1:A:455:MET:HB3	2.01	0.41
1:A:683:ASN:HA	1:A:709:TRP:CD1	2.56	0.41
1:A:786:PHE:CE2	1:A:790:ILE:HD11	2.56	0.41
1:B:312:LYS:O	1:B:316:GLU:HG3	2.20	0.41
1:B:518:ARG:HH22	1:B:521:ASP:CG	2.22	0.41
1:B:545:ILE:HD13	1:B:903:LEU:HD23	2.02	0.41
1:B:330:THR:HB	1:B:333:LYS:CG	2.50	0.41
1:B:342:ILE:HG23	1:B:352:ALA:HB2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:663:MET:HG3	1:B:904:ILE:HD11	2.02	0.41
1:A:356:LEU:HB3	1:A:363:PRO:HG3	2.03	0.41
1:B:98:ASN:CB	1:B:103:GLN:HB2	2.47	0.41
1:B:193:ILE:HD13	1:B:201:ALA:CB	2.50	0.41
1:B:361:VAL:O	1:B:363:PRO:HD3	2.21	0.41
1:B:786:PHE:HZ	1:B:804:LEU:CD2	2.34	0.41
1:B:94:ARG:O	1:B:107:MET:HA	2.21	0.41
1:B:83:LEU:HD21	1:B:134:LEU:HD13	2.02	0.41
1:B:83:LEU:HB2	1:B:152:PHE:CD1	2.56	0.41
1:B:400:ASN:HD22	1:B:400:ASN:HA	1.70	0.41
1:A:531:LEU:HD21	1:A:582:ILE:HD11	2.02	0.40
1:A:551:ARG:HG3	1:A:551:ARG:NH1	2.35	0.40
1:B:490:ARG:O	1:B:494:GLU:HG2	2.20	0.40
1:A:243:ARG:NH1	1:A:244:HIS:HE1	2.18	0.40
1:B:709:TRP:C	1:B:709:TRP:CD1	2.94	0.40
1:B:806:GLN:HB3	1:B:806:GLN:HE21	1.56	0.40
1:A:338:ASP:O	1:A:342:ILE:HD13	2.21	0.40
1:A:637:LEU:HD23	1:A:651:VAL:HG21	2.01	0.40
1:B:531:LEU:HD13	1:B:598:LEU:HD11	2.03	0.40
1:A:46:GLU:HG3	1:A:264:PHE:HE1	1.86	0.40
1:A:520:PRO:HD3	1:A:663:MET:CE	2.52	0.40
1:A:152:PHE:O	1:A:206:VAL:HA	2.22	0.40
1:B:767:LEU:CG	1:B:818:LEU:HD23	2.47	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	897/917 (98%)	840 (94%)	55 (6%)	2 (0%)	51 71
1	B	897/917 (98%)	845 (94%)	49 (6%)	3 (0%)	44 64

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1794/1834 (98%)	1685 (94%)	104 (6%)	5 (0%)	44 64

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	104	ASN
1	B	102	ASN
1	A	243	ARG
1	A	450	GLY
1	B	436	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	774/788 (98%)	749 (97%)	25 (3%)	44 69
1	B	774/788 (98%)	756 (98%)	18 (2%)	56 79
All	All	1548/1576 (98%)	1505 (97%)	43 (3%)	49 74

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	ASP
1	A	21	LYS
1	A	27	TYR
1	A	96	GLN
1	A	102	ASN
1	A	159	GLN
1	A	196	ARG
1	A	198	ASP
1	A	199	TYR
1	A	202	ASN
1	A	254	ARG
1	A	261	TRP
1	A	320	PHE

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Mol	Chain	Res	Type
1	A	400	ASN
1	A	407	ARG
1	A	408	THR
1	A	481	LYS
1	A	502	GLN
1	A	531	LEU
1	A	565	GLU
1	A	709	TRP
1	A	810	ASN
1	A	843	ASP
1	A	857	THR
1	A	913	THR
1	B	66	THR
1	B	99	HIS
1	B	202	ASN
1	B	261	TRP
1	B	320	PHE
1	B	321	GLU
1	B	380	PHE
1	B	407	ARG
1	B	424	SER
1	B	437	ASP
1	B	463	LEU
1	B	709	TRP
1	B	769	ARG
1	B	798	LEU
1	B	805	GLN
1	B	806	GLN
1	B	810	ASN
1	B	843	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (39) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	102	ASN
1	A	104	ASN
1	A	125	GLN
1	A	159	GLN
1	A	202	ASN
1	A	244	HIS
1	A	345	ASN
1	A	351	ASN

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Mol	Chain	Res	Type
1	A	384	ASN
1	A	400	ASN
1	A	420	HIS
1	A	466	GLN
1	A	506	ASN
1	A	577	HIS
1	A	700	GLN
1	A	771	GLN
1	A	805	GLN
1	A	806	GLN
1	A	810	ASN
1	A	887	ASN
1	B	98	ASN
1	B	104	ASN
1	B	202	ASN
1	B	244	HIS
1	B	345	ASN
1	B	384	ASN
1	B	400	ASN
1	B	420	HIS
1	B	466	GLN
1	B	502	GLN
1	B	506	ASN
1	B	557	ASN
1	B	702	GLN
1	B	771	GLN
1	B	799	GLN
1	B	806	GLN
1	B	810	ASN
1	B	832	GLN
1	B	887	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 14 ligands modelled in this entry, 4 are monoatomic - leaving 10 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$
2	BGC	A	1001	-	12,12,12	0.51	0	17,17,17	1.50	2 (11%)
3	BG6	A	1002	-	16,16,16	1.10	1 (6%)	24,24,24	0.82	1 (4%)
2	BGC	A	1003	-	12,12,12	0.37	0	17,17,17	1.26	2 (11%)
3	BG6	A	1004	-	16,16,16	1.01	1 (6%)	24,24,24	0.97	1 (4%)
5	CIT	A	1007	-	3,12,12	1.22	0	3,17,17	1.60	1 (33%)
2	BGC	B	1001	-	12,12,12	0.37	0	17,17,17	1.42	2 (11%)
3	BG6	B	1002	-	16,16,16	1.02	1 (6%)	24,24,24	0.73	1 (4%)
2	BGC	B	1003	-	12,12,12	0.48	0	17,17,17	1.35	3 (17%)
3	BG6	B	1004	-	16,16,16	1.06	1 (6%)	24,24,24	0.83	1 (4%)
5	CIT	B	1007	-	3,12,12	1.28	0	3,17,17	1.74	1 (33%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BGC	A	1001	-	-	0/2/22/22	0/1/1/1
3	BG6	A	1002	-	-	0/6/26/26	0/1/1/1
2	BGC	A	1003	-	-	0/2/22/22	0/1/1/1
3	BG6	A	1004	-	-	0/6/26/26	0/1/1/1
5	CIT	A	1007	-	-	0/6/16/16	0/0/0/0
2	BGC	B	1001	-	-	0/2/22/22	0/1/1/1
3	BG6	B	1002	-	-	0/6/26/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BGC	B	1003	-	-	0/2/22/22	0/1/1/1
3	BG6	B	1004	-	-	0/6/26/26	0/1/1/1
5	CIT	B	1007	-	-	0/6/16/16	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1004	BG6	P-O3P	2.96	1.60	1.50
3	B	1004	BG6	P-O3P	3.09	1.61	1.50
3	B	1002	BG6	P-O3P	3.17	1.61	1.50
3	A	1002	BG6	P-O3P	3.36	1.62	1.50

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	BGC	C1-O5-C5	-4.43	105.41	113.39
2	B	1001	BGC	C1-O5-C5	-3.73	106.67	113.39
2	B	1001	BGC	O5-C1-C2	-3.69	103.92	110.04
2	A	1003	BGC	O5-C1-C2	-3.31	104.55	110.04
2	B	1003	BGC	C1-O5-C5	-3.25	107.53	113.39
2	A	1001	BGC	O5-C1-C2	-3.22	104.69	110.04
2	B	1003	BGC	O5-C1-C2	-3.19	104.74	110.04
2	A	1003	BGC	C1-O5-C5	-2.89	108.17	113.39
5	B	1007	CIT	C3-C2-C1	-2.63	110.84	114.95
5	A	1007	CIT	C3-C4-C5	-2.37	111.24	114.95
2	B	1003	BGC	C1-C2-C3	-2.04	106.96	110.65
3	B	1002	BG6	O1P-P-O6	2.14	112.44	106.73
3	A	1002	BG6	O1P-P-O6	2.60	113.65	106.73
3	B	1004	BG6	O1P-P-O6	2.77	114.09	106.73
3	A	1004	BG6	O1P-P-O6	3.03	114.79	106.73

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1002	BG6	2	0
2	A	1003	BGC	2	0
5	A	1007	CIT	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1004	BG6	1	0
5	B	1007	CIT	5	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	899/917 (98%)	0.21	32 (3%) 43 45	31, 59, 85, 113	0
1	B	899/917 (98%)	0.21	29 (3%) 48 50	32, 58, 85, 112	0
All	All	1798/1834 (98%)	0.21	61 (3%) 46 47	31, 59, 85, 113	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	22	ILE	5.4
1	A	16	ASP	4.6
1	B	16	ASP	4.3
1	B	19	VAL	4.3
1	A	564	ILE	4.0
1	A	21	LYS	3.7
1	A	794	ARG	3.7
1	B	18	GLN	3.6
1	A	17	ASP	3.5
1	A	786	PHE	3.4
1	B	102	ASN	3.4
1	B	100	GLU	3.4
1	A	31	LEU	3.3
1	B	794	ARG	3.2
1	B	913	THR	3.1
1	B	786	PHE	3.1
1	A	443	LEU	2.8
1	A	310	LEU	2.8
1	A	350	HIS	2.8
1	A	229	ILE	2.8
1	B	17	ASP	2.7
1	B	359	LEU	2.7
1	B	103	GLN	2.7
1	B	237	CYS	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	101	LYS	2.6
1	A	26	LEU	2.6
1	B	350	HIS	2.6
1	A	436	PRO	2.5
1	B	346	LYS	2.5
1	B	798	LEU	2.5
1	A	417	TYR	2.4
1	A	911	LEU	2.4
1	A	30	ARG	2.4
1	A	142	LYS	2.3
1	A	199	TYR	2.3
1	B	324	ILE	2.3
1	A	362	GLU	2.3
1	A	358	ARG	2.2
1	A	769	ARG	2.2
1	B	235	ASN	2.2
1	A	19	VAL	2.2
1	B	318	LEU	2.2
1	B	549	LYS	2.2
1	B	262	GLY	2.2
1	B	406	LEU	2.2
1	B	548	GLY	2.2
1	B	518	ARG	2.2
1	B	234	THR	2.2
1	B	258	ASN	2.2
1	B	550	LYS	2.1
1	B	263	ALA	2.1
1	B	227	GLY	2.1
1	A	412	VAL	2.1
1	A	204	VAL	2.1
1	A	320	PHE	2.0
1	A	410	VAL	2.0
1	A	359	LEU	2.0
1	A	414	GLY	2.0
1	A	424	SER	2.0
1	A	444	LEU	2.0
1	A	549	LYS	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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
5	CIT	B	1007	13/13	0.56	0.27	5.76	106,108,108,108	0
2	BGC	B	1001	12/12	0.94	0.33	2.43	48,49,50,51	0
2	BGC	A	1001	12/12	0.89	0.27	1.49	50,52,53,55	0
2	BGC	A	1003	12/12	0.96	0.21	1.35	36,38,38,41	0
3	BG6	A	1002	16/16	0.93	0.23	0.45	75,78,80,80	0
3	BG6	A	1004	16/16	0.98	0.18	0.35	36,41,42,43	0
2	BGC	B	1003	12/12	0.93	0.18	0.33	38,39,40,42	0
3	BG6	B	1004	16/16	0.96	0.18	0.24	36,43,44,45	0
3	BG6	B	1002	16/16	0.95	0.21	0.09	71,74,75,76	0
4	NA	B	1006	1/1	0.92	0.07	-1.83	52,52,52,52	0
4	NA	B	1005	1/1	0.90	0.09	-1.84	65,65,65,65	0
4	NA	A	1006	1/1	0.91	0.05	-1.96	53,53,53,53	0
4	NA	A	1005	1/1	0.89	0.08	-2.63	63,63,63,63	0
5	CIT	A	1007	13/13	0.72	0.36	-	110,110,111,111	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.