



Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 02:33 AM GMT

PDB ID : 2C2L
Title : CRYSTAL STRUCTURE OF THE CHIP U-BOX E3 UBIQUITIN LIGASE
Authors : Zhang, M.; Roe, S.M.; Pearl, L.H.
Deposited on : 2005-09-29
Resolution : 3.30 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

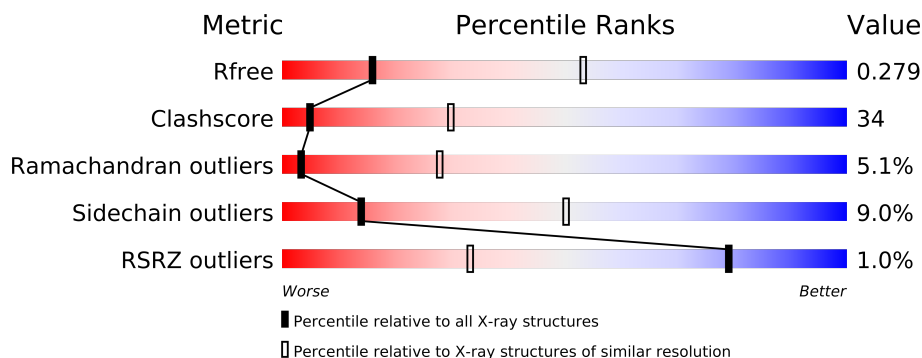
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.30 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}	66092	1341 (3.40-3.20)
Clashscore	79885	1696 (3.40-3.20)
Ramachandran outliers	78287	1664 (3.40-3.20)
Sidechain outliers	78261	1662 (3.40-3.20)
RSRZ outliers	66119	1342 (3.40-3.20)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	281	
1	B	281	
1	C	281	
1	D	281	
2	E	9	
2	F	9	
2	G	9	
2	H	9	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	SO4	A	1307	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9521 atoms, of which 0 are hydrogen and 0 are deuterium.

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 CARBOXY TERMINUS OF HSP70-INTERACTING PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	281	Total	C	N	O	S	0	0	0
			2294	1424	417	439	14			
1	B	281	Total	C	N	O	S	0	0	0
			2294	1424	417	439	14			
1	C	281	Total	C	N	O	S	0	0	0
			2294	1424	417	439	14			
1	D	281	Total	C	N	O	S	0	0	0
			2294	1424	417	439	14			

- Molecule 2 is a protein called HSP90.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	9	Total	C	N	O	S	0	0	0
			74	41	12	20	1			
2	F	9	Total	C	N	O	S	0	0	0
			74	41	12	20	1			
2	G	9	Total	C	N	O	S	0	0	0
			74	41	12	20	1			
2	H	9	Total	C	N	O	S	0	0	0
			74	41	12	20	1			

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	4	Total	Ni	0	0
			4	4		

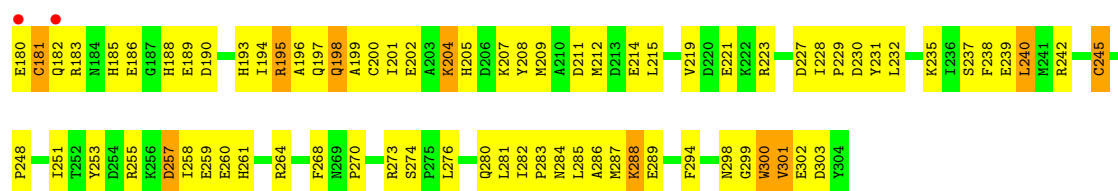
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	4	Total	O	0	0
			4	4		

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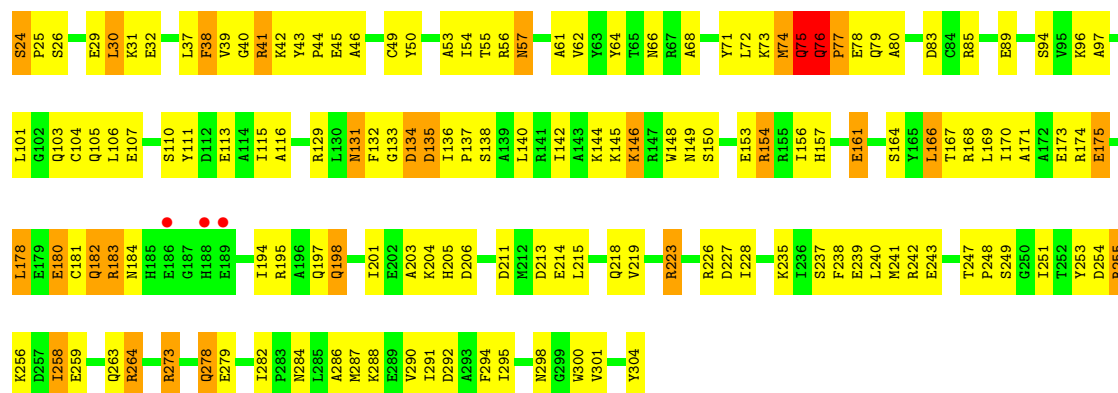
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	3	Total 3	O 3	0	0
5	D	2	Total 2	O 2	0	0
5	G	1	Total 1	O 1	0	0



• Molecule 1: CARBOXY TERMINUS OF HSP70-INTERACTING PROTEIN

Chain D:



• Molecule 2: HSP90

Chain E:



• Molecule 2: HSP90

Chain F:



• Molecule 2: HSP90

Chain G:



• Molecule 2: HSP90

Chain H:



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	76.04Å 204.41Å 144.71Å 90.00° 90.75° 90.00°	Depositor
Resolution (Å)	40.00 – 3.30 30.05 – 3.31	Depositor EDS
% Data completeness (in resolution range)	21.5 (40.00-3.30) 94.0 (30.05-3.31)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 3.31Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.247 , 0.286 0.242 , 0.279	Depositor DCC
R_{free} test set	1606 reflections (5.03%)	DCC
Wilson B-factor (Å ²)	78.0	Xtriage
Anisotropy	0.815	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 39.6	EDS
Estimated twinning fraction	0.327 for h,-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 31915 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9521	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.94% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: NI, SO4

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	0.50	1/2337 (0.0%)	0.60	0/3144
1	B	0.50	0/2337	0.62	0/3144
1	C	0.50	1/2337 (0.0%)	0.60	0/3144
1	D	0.51	0/2337	0.62	0/3144
2	E	0.60	0/73	0.61	0/95
2	F	0.59	0/73	0.58	0/95
2	G	0.60	0/73	0.60	0/95
2	H	0.62	0/73	0.58	0/95
All	All	0.51	2/9640 (0.0%)	0.61	0/12956

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	245	CYS	CB-SG	-6.38	1.71	1.82
1	C	245	CYS	CB-SG	-5.95	1.72	1.81

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2294	0	2234	147	0
1	B	2294	0	2234	183	0
1	C	2294	0	2234	151	0
1	D	2294	0	2234	175	0
2	E	74	0	62	10	0
2	F	74	0	62	15	0
2	G	74	0	62	10	0
2	H	74	0	62	12	0
3	A	15	0	0	2	0
3	B	10	0	0	0	0
3	C	5	0	0	0	0
3	D	5	0	0	0	0
4	B	4	0	0	0	0
5	A	4	0	0	0	0
5	C	3	0	0	0	0
5	D	2	0	0	0	0
5	G	1	0	0	0	0
All	All	9521	0	9184	634	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 34.

All (634) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:74:MET:HG3	1:D:75:GLN:H	1.12	1.05
1:B:74:MET:HG3	1:B:75:GLN:H	1.11	1.05
1:D:96:LYS:HG3	2:H:505:MET:HE2	1.40	1.02
1:D:71:TYR:HB3	1:D:76:GLN:HB3	1.42	1.01
1:B:71:TYR:HB3	1:B:76:GLN:HB3	1.43	0.99
1:B:96:LYS:HG3	2:F:505:MET:HE2	1.43	0.99
1:C:194:ILE:HA	1:C:197:GLN:HB2	1.43	0.96
1:D:138:SER:O	1:D:142:ILE:HD12	1.66	0.96
1:A:194:ILE:HA	1:A:197:GLN:HB2	1.44	0.95
1:B:138:SER:O	1:B:142:ILE:HD12	1.67	0.95
1:A:182:GLN:HE21	1:A:198:GLN:HG3	1.36	0.91
1:B:56:ARG:HH22	1:C:56:ARG:HH21	0.96	0.90
1:A:151:ILE:H	1:A:151:ILE:HD13	1.36	0.88
1:C:182:GLN:HE21	1:C:198:GLN:HG3	1.39	0.88
1:B:258:ILE:HD12	1:B:259:GLU:H	1.37	0.87
1:A:221:GLU:HB3	1:A:223:ARG:HG2	1.57	0.87
1:C:151:ILE:HD13	1:C:151:ILE:H	1.37	0.86
1:C:221:GLU:HB3	1:C:223:ARG:HG2	1.58	0.86
1:D:258:ILE:HD12	1:D:259:GLU:H	1.38	0.86

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:74:MET:HG3	1:B:75:GLN:N	1.91	0.85
1:D:74:MET:HG3	1:D:75:GLN:N	1.92	0.84
1:C:160:SER:HB3	1:C:163:HIS:HB2	1.57	0.84
1:A:160:SER:HB3	1:A:163:HIS:HB2	1.58	0.83
1:B:56:ARG:NH2	1:C:56:ARG:HH21	1.76	0.83
1:C:164:SER:O	1:C:167:THR:HG22	1.79	0.82
1:B:75:GLN:C	1:B:77:PRO:HD2	1.98	0.82
1:D:75:GLN:C	1:D:77:PRO:HD2	1.98	0.82
1:A:164:SER:O	1:A:167:THR:HG22	1.79	0.82
1:D:26:SER:HA	1:D:56:ARG:HH12	1.45	0.81
1:B:26:SER:HA	1:B:56:ARG:HH12	1.45	0.81
1:B:131:ASN:HD22	1:B:131:ASN:H	1.28	0.81
1:B:258:ILE:HD12	1:B:259:GLU:N	1.96	0.80
1:D:258:ILE:HD12	1:D:259:GLU:N	1.96	0.80
1:D:37:LEU:HB2	1:D:46:ALA:HB2	1.63	0.79
1:D:71:TYR:HB3	1:D:76:GLN:CB	2.12	0.79
1:D:131:ASN:HD22	1:D:131:ASN:H	1.29	0.79
1:B:71:TYR:HB3	1:B:76:GLN:CB	2.12	0.79
1:B:174:ARG:HG2	1:B:175:GLU:OE1	1.82	0.78
1:D:174:ARG:HG2	1:D:175:GLU:OE1	1.82	0.78
1:B:37:LEU:HB2	1:B:46:ALA:HB2	1.64	0.78
1:B:181:CYS:HA	1:B:184:ASN:HD22	1.49	0.78
1:D:181:CYS:HA	1:D:184:ASN:HD22	1.49	0.78
1:D:149:ASN:O	1:D:153:GLU:HG3	1.84	0.77
1:C:248:PRO:HA	1:D:282:ILE:HD13	1.66	0.77
1:B:149:ASN:O	1:B:153:GLU:HG3	1.85	0.77
1:D:43:TYR:N	1:D:44:PRO:HD2	1.99	0.77
1:B:93:GLN:HE21	1:C:93:GLN:HG3	1.50	0.77
1:B:43:TYR:N	1:B:44:PRO:HD2	1.99	0.76
1:C:301:VAL:C	1:C:303:ASP:H	1.87	0.76
1:C:255:ARG:O	1:C:259:GLU:HB2	1.86	0.76
1:A:301:VAL:C	1:A:303:ASP:H	1.88	0.75
1:A:284:ASN:HD21	1:B:284:ASN:HD21	1.32	0.74
1:A:255:ARG:O	1:A:259:GLU:HB2	1.88	0.74
1:B:72:LEU:HD23	1:B:76:GLN:HG2	1.71	0.73
1:D:66:ASN:CG	2:H:508:VAL:HG13	2.09	0.73
1:B:76:GLN:N	1:B:77:PRO:HD2	2.04	0.73
1:B:56:ARG:HH22	1:C:56:ARG:NH2	1.80	0.72
1:D:72:LEU:HD23	1:D:76:GLN:HG2	1.71	0.72
1:C:78:GLU:HB2	1:C:79:GLN:NE2	2.04	0.72
1:D:76:GLN:N	1:D:77:PRO:HD2	2.05	0.72
1:A:78:GLU:HB2	1:A:79:GLN:NE2	2.05	0.71

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:182:GLN:NE2	1:A:198:GLN:HG3	2.04	0.71
1:B:194:ILE:HA	1:B:197:GLN:HB3	1.73	0.70
1:A:248:PRO:HA	1:B:282:ILE:HD13	1.73	0.70
1:D:194:ILE:HA	1:D:197:GLN:HB3	1.74	0.70
1:D:194:ILE:O	1:D:198:GLN:HB2	1.91	0.69
1:C:66:ASN:CG	2:G:508:VAL:HG13	2.13	0.69
1:D:50:TYR:OH	2:H:508:VAL:HG11	1.92	0.68
1:C:284:ASN:HD21	1:D:284:ASN:HD21	1.38	0.68
1:A:66:ASN:CG	2:E:508:VAL:HG13	2.13	0.68
1:B:182:GLN:HG3	1:B:198:GLN:NE2	2.09	0.68
1:D:37:LEU:CB	1:D:46:ALA:HB2	2.24	0.68
1:B:194:ILE:O	1:B:198:GLN:HB2	1.92	0.68
1:A:284:ASN:HD21	1:B:284:ASN:ND2	1.92	0.68
1:B:110:SER:HB3	1:B:113:GLU:OE1	1.94	0.68
1:C:182:GLN:NE2	1:C:198:GLN:HG3	2.08	0.68
1:C:212:MET:O	1:C:215:LEU:HB2	1.93	0.68
1:B:264:ARG:HD3	1:B:264:ARG:O	1.94	0.68
1:C:134:ASP:O	1:C:138:SER:HB3	1.94	0.67
1:A:134:ASP:O	1:A:138:SER:HB3	1.94	0.67
1:D:110:SER:HB3	1:D:113:GLU:OE1	1.94	0.67
1:B:103:GLN:OE1	1:B:103:GLN:HA	1.93	0.67
1:B:37:LEU:CB	1:B:46:ALA:HB2	2.25	0.67
1:D:145:LYS:HD3	1:D:304:TYR:CD2	2.29	0.67
1:C:251:ILE:N	1:C:251:ILE:HD12	2.09	0.67
1:B:94:SER:HB3	1:B:97:ALA:HB3	1.76	0.67
1:A:284:ASN:ND2	1:B:284:ASN:HD21	1.91	0.67
1:D:264:ARG:HD3	1:D:264:ARG:O	1.94	0.67
1:B:145:LYS:HD3	1:B:304:TYR:CD2	2.30	0.67
1:A:212:MET:O	1:A:215:LEU:HB2	1.95	0.67
1:D:94:SER:HB3	1:D:97:ALA:HB3	1.77	0.66
1:A:251:ILE:HD12	1:A:251:ILE:N	2.10	0.66
1:B:255:ARG:O	1:B:259:GLU:HB2	1.95	0.66
1:D:182:GLN:HG3	1:D:198:GLN:NE2	2.11	0.66
1:D:103:GLN:HA	1:D:103:GLN:OE1	1.95	0.65
1:C:103:GLN:O	1:C:106:LEU:HB2	1.97	0.65
1:A:182:GLN:HG2	1:A:198:GLN:HE21	1.61	0.65
2:E:508:VAL:O	2:E:508:VAL:HG12	1.96	0.65
1:A:103:GLN:O	1:A:106:LEU:HB2	1.97	0.65
1:D:255:ARG:O	1:D:259:GLU:HB2	1.96	0.65
1:B:182:GLN:HG3	1:B:198:GLN:HE22	1.59	0.64
2:G:508:VAL:O	2:G:508:VAL:HG12	1.96	0.64
1:B:50:TYR:OH	2:F:508:VAL:HG11	1.97	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:508:VAL:HG12	2:F:508:VAL:O	1.97	0.64
2:H:508:VAL:O	2:H:508:VAL:HG12	1.97	0.64
1:B:66:ASN:CG	2:F:508:VAL:HG13	2.18	0.64
1:C:182:GLN:HG2	1:C:198:GLN:HE21	1.62	0.63
1:B:154:ARG:HD3	1:B:154:ARG:O	1.99	0.63
1:A:148:TRP:HA	1:A:151:ILE:HD11	1.80	0.63
1:B:284:ASN:OD1	1:B:287:MET:HB2	1.98	0.63
1:A:201:ILE:HD12	1:A:204:LYS:HB3	1.80	0.63
1:C:183:ARG:HG3	1:C:186:GLU:HG3	1.81	0.63
1:D:154:ARG:O	1:D:154:ARG:HD3	1.99	0.63
1:C:78:GLU:HB2	1:C:79:GLN:HE21	1.62	0.63
1:A:78:GLU:HB2	1:A:79:GLN:HE21	1.62	0.63
1:A:183:ARG:HG3	1:A:186:GLU:HG3	1.81	0.62
1:D:131:ASN:HB2	2:H:502:THR:HG23	1.81	0.62
1:B:131:ASN:HD22	1:B:131:ASN:N	1.95	0.62
1:D:182:GLN:HG3	1:D:198:GLN:HE22	1.62	0.62
1:C:284:ASN:ND2	1:D:284:ASN:HD21	1.96	0.62
1:A:85:ARG:HH11	1:A:85:ARG:HG3	1.64	0.62
1:C:148:TRP:HA	1:C:151:ILE:HD11	1.82	0.62
1:D:85:ARG:HH11	1:D:85:ARG:HG3	1.64	0.62
1:D:284:ASN:OD1	1:D:287:MET:HB2	2.00	0.61
1:C:185:HIS:HB3	1:C:188:HIS:HB2	1.82	0.61
1:C:85:ARG:HG3	1:C:85:ARG:HH11	1.65	0.61
1:A:189:GLU:HG2	1:A:193:HIS:ND1	2.16	0.61
1:A:185:HIS:HB3	1:A:188:HIS:HB2	1.82	0.61
1:C:156:ILE:N	1:C:156:ILE:HD12	2.16	0.61
1:A:166:LEU:HD23	1:A:169:LEU:HD12	1.83	0.61
1:B:85:ARG:HG3	1:B:85:ARG:HH11	1.65	0.61
1:A:156:ILE:N	1:A:156:ILE:HD12	2.16	0.60
1:C:151:ILE:HD13	1:C:151:ILE:N	2.15	0.60
1:B:291:ILE:O	1:B:295:ILE:HD12	2.00	0.60
1:C:201:ILE:HD12	1:C:204:LYS:HB3	1.83	0.60
1:C:111:TYR:O	1:C:115:ILE:HG23	2.01	0.60
1:C:142:ILE:O	1:C:146:LYS:HG3	2.02	0.60
1:C:96:LYS:HG3	2:G:505:MET:HE1	1.84	0.60
1:C:189:GLU:HG2	1:C:193:HIS:ND1	2.17	0.60
1:D:71:TYR:CE2	1:D:79:GLN:HB3	2.37	0.60
1:B:71:TYR:CE2	1:B:79:GLN:HB3	2.37	0.60
1:B:111:TYR:HE2	1:B:146:LYS:HB3	1.67	0.60
1:D:74:MET:CG	1:D:75:GLN:H	1.99	0.59
1:C:166:LEU:HD23	1:C:169:LEU:HD12	1.84	0.59
1:C:284:ASN:HD21	1:D:284:ASN:ND2	2.00	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:131:ASN:ND2	2:G:502:THR:HG23	2.17	0.59
1:C:228:ILE:HD12	1:C:229:PRO:O	2.02	0.59
1:D:111:TYR:HE2	1:D:146:LYS:HB3	1.68	0.59
1:A:142:ILE:O	1:A:146:LYS:HG3	2.03	0.59
1:B:74:MET:CG	1:B:75:GLN:H	1.98	0.59
1:A:165:TYR:O	1:A:168:ARG:HG2	2.02	0.59
1:D:164:SER:O	1:D:167:THR:HG22	2.03	0.59
1:A:228:ILE:HD12	1:A:229:PRO:O	2.02	0.59
1:B:182:GLN:HE21	1:B:198:GLN:NE2	2.01	0.58
1:B:145:LYS:CE	1:B:304:TYR:HB3	2.33	0.58
1:B:54:ILE:HD13	1:B:64:TYR:CE1	2.38	0.58
1:B:134:ASP:HA	1:B:137:PRO:HG2	1.86	0.58
1:A:111:TYR:O	1:A:115:ILE:HG23	2.03	0.58
1:C:212:MET:HA	1:C:215:LEU:HD12	1.85	0.58
1:A:212:MET:HA	1:A:215:LEU:HD12	1.85	0.58
1:D:211:ASP:O	1:D:214:GLU:HB3	2.03	0.58
1:B:164:SER:O	1:B:167:THR:HG22	2.04	0.58
1:C:154:ARG:HH11	1:C:154:ARG:HG2	1.69	0.58
1:D:57:ASN:HD22	1:D:57:ASN:C	2.07	0.58
1:D:134:ASP:HA	1:D:137:PRO:HG2	1.86	0.58
1:B:72:LEU:HA	1:B:76:GLN:HG3	1.85	0.57
1:D:145:LYS:HD3	1:D:304:TYR:HD2	1.67	0.57
1:B:57:ASN:C	1:B:57:ASN:HD22	2.07	0.57
1:B:132:PHE:HB3	1:B:135:ASP:HB2	1.86	0.57
1:D:145:LYS:CE	1:D:304:TYR:HB3	2.34	0.57
1:C:98:HIS:HB3	1:C:121:ALA:HB2	1.87	0.57
1:D:54:ILE:HD13	1:D:64:TYR:CE1	2.39	0.57
1:C:165:TYR:O	1:C:168:ARG:HG2	2.04	0.57
1:C:160:SER:HB3	1:C:163:HIS:CB	2.33	0.57
1:C:174:ARG:HH21	1:C:205:HIS:CB	2.17	0.57
1:A:154:ARG:HH11	1:A:154:ARG:HG2	1.70	0.57
1:B:237:SER:O	1:B:238:PHE:HB2	2.04	0.57
1:A:160:SER:HB3	1:A:163:HIS:CB	2.33	0.57
1:B:145:LYS:HD3	1:B:304:TYR:HD2	1.68	0.57
1:B:211:ASP:O	1:B:214:GLU:HB3	2.04	0.57
1:A:174:ARG:HH21	1:A:205:HIS:CB	2.18	0.57
1:D:291:ILE:O	1:D:295:ILE:HD12	2.04	0.57
1:B:140:LEU:O	1:B:144:LYS:HG2	2.05	0.57
1:D:140:LEU:O	1:D:144:LYS:HG2	2.05	0.57
1:A:98:HIS:HB3	1:A:121:ALA:HB2	1.87	0.56
1:A:189:GLU:HG2	1:A:193:HIS:CE1	2.41	0.56
1:D:72:LEU:HA	1:D:76:GLN:HG3	1.86	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:131:ASN:CG	2:H:502:THR:HG23	2.26	0.56
1:A:196:ALA:O	1:A:200:CYS:HB2	2.05	0.56
1:C:299:GLY:O	1:C:300:TRP:O	2.23	0.56
1:C:28:GLN:O	1:C:32:GLU:HG3	2.04	0.56
1:D:132:PHE:HB3	1:D:135:ASP:HB2	1.87	0.56
1:A:151:ILE:H	1:A:151:ILE:CD1	2.12	0.56
1:D:29:GLU:O	1:D:31:LYS:N	2.39	0.56
1:C:139:ALA:HA	1:C:142:ILE:HD12	1.88	0.56
1:A:28:GLN:O	1:A:32:GLU:HG3	2.04	0.56
1:B:85:ARG:O	1:B:89:GLU:HG3	2.05	0.56
1:D:66:ASN:OD1	2:H:508:VAL:HG13	2.06	0.56
1:A:139:ALA:HA	1:A:142:ILE:HD12	1.88	0.56
1:D:237:SER:O	1:D:238:PHE:HB2	2.05	0.56
1:B:180:GLU:HG3	1:B:181:CYS:N	2.21	0.56
1:D:180:GLU:HG3	1:D:181:CYS:N	2.21	0.56
1:A:282:ILE:HD12	1:B:282:ILE:HD12	1.88	0.55
1:D:85:ARG:O	1:D:89:GLU:HG3	2.06	0.55
1:B:298:ASN:O	1:B:301:VAL:HG23	2.06	0.55
1:A:174:ARG:HH21	1:A:205:HIS:HB2	1.71	0.55
1:A:294:PHE:O	1:A:298:ASN:ND2	2.39	0.55
1:A:131:ASN:ND2	2:E:502:THR:HG23	2.20	0.55
1:D:43:TYR:N	1:D:44:PRO:CD	2.67	0.55
1:A:36:ARG:HD2	3:A:1307:SO4:O1	2.06	0.55
1:A:164:SER:C	1:A:167:THR:HG22	2.26	0.55
1:C:255:ARG:HA	1:C:258:ILE:HD11	1.89	0.55
1:B:60:VAL:HG22	1:C:59:LEU:O	2.07	0.55
1:B:43:TYR:N	1:B:44:PRO:CD	2.67	0.55
1:C:294:PHE:O	1:C:298:ASN:ND2	2.39	0.55
1:C:189:GLU:HG2	1:C:193:HIS:CE1	2.42	0.55
1:C:164:SER:C	1:C:167:THR:HG22	2.27	0.55
1:A:255:ARG:HA	1:A:258:ILE:HD11	1.89	0.55
1:C:174:ARG:HH21	1:C:205:HIS:HB2	1.72	0.55
1:C:301:VAL:O	1:C:302:GLU:HB3	2.06	0.54
1:A:176:ARG:HH12	1:B:173:GLU:HG2	1.72	0.54
1:A:42:LYS:HE2	1:C:283:PRO:CD	2.37	0.54
1:B:131:ASN:HB2	2:F:502:THR:HG23	1.89	0.54
1:D:182:GLN:HE21	1:D:198:GLN:NE2	2.06	0.54
1:A:301:VAL:O	1:A:302:GLU:HB3	2.06	0.54
1:A:219:VAL:HG13	1:A:219:VAL:O	2.07	0.54
1:D:131:ASN:CB	2:H:502:THR:HG23	2.37	0.54
1:A:301:VAL:C	1:A:303:ASP:N	2.57	0.54
1:C:196:ALA:O	1:C:200:CYS:HB2	2.08	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:151:ILE:HD13	1:A:151:ILE:N	2.14	0.54
1:D:201:ILE:HD12	1:D:204:LYS:HZ2	1.72	0.54
2:H:505:MET:C	2:H:507:GLU:H	2.11	0.54
1:C:301:VAL:C	1:C:303:ASP:N	2.56	0.54
1:C:219:VAL:O	1:C:219:VAL:HG22	2.07	0.54
1:B:181:CYS:O	1:B:184:ASN:HB2	2.07	0.54
1:B:93:GLN:NE2	1:C:93:GLN:HG3	2.21	0.54
1:B:69:LEU:HD23	2:F:508:VAL:CG2	2.37	0.54
1:D:298:ASN:O	1:D:301:VAL:HG23	2.07	0.54
1:C:151:ILE:H	1:C:151:ILE:CD1	2.13	0.54
1:A:189:GLU:HB3	1:A:194:ILE:HG23	1.89	0.54
1:D:42:LYS:C	1:D:44:PRO:HD2	2.28	0.54
1:D:181:CYS:O	1:D:184:ASN:HB2	2.08	0.54
1:B:111:TYR:CE2	1:B:146:LYS:HB3	2.43	0.54
1:A:24:SER:N	1:A:25:PRO:HD2	2.23	0.54
2:E:505:MET:C	2:E:507:GLU:H	2.11	0.54
1:D:68:ALA:HB2	1:D:83:ASP:HB3	1.90	0.54
1:B:26:SER:HB3	1:B:29:GLU:CD	2.28	0.53
1:B:42:LYS:C	1:B:44:PRO:HD2	2.28	0.53
1:C:25:PRO:CG	1:C:30:LEU:HD21	2.38	0.53
1:B:273:ARG:HH11	1:B:273:ARG:HG2	1.73	0.53
1:B:136:ILE:N	1:B:137:PRO:HD2	2.23	0.53
1:C:219:VAL:O	1:C:219:VAL:HG13	2.09	0.53
1:D:29:GLU:C	1:D:31:LYS:H	2.12	0.53
1:A:299:GLY:O	1:A:300:TRP:O	2.26	0.53
1:C:212:MET:O	1:C:215:LEU:N	2.38	0.53
1:A:25:PRO:CG	1:A:30:LEU:HD21	2.38	0.53
1:D:76:GLN:H	1:D:76:GLN:CD	2.10	0.53
1:D:111:TYR:CE2	1:D:146:LYS:HB3	2.44	0.53
1:C:24:SER:N	1:C:25:PRO:HD2	2.24	0.53
2:F:505:MET:C	2:F:507:GLU:H	2.12	0.53
2:G:505:MET:C	2:G:507:GLU:H	2.12	0.53
1:B:68:ALA:HB2	1:B:83:ASP:HB3	1.91	0.53
1:A:176:ARG:NH1	1:B:173:GLU:HG2	2.24	0.52
1:B:29:GLU:O	1:B:31:LYS:N	2.42	0.52
1:B:241:MET:HE1	1:B:253:TYR:HA	1.90	0.52
1:D:241:MET:HE1	1:D:253:TYR:HA	1.90	0.52
1:B:178:LEU:O	1:B:182:GLN:HB3	2.09	0.52
1:A:96:LYS:HG3	2:E:505:MET:HE1	1.91	0.52
1:B:24:SER:HB2	1:B:25:PRO:HD3	1.90	0.52
1:C:167:THR:HG23	1:C:168:ARG:N	2.25	0.52
1:B:76:GLN:H	1:B:76:GLN:CD	2.11	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:136:ILE:N	1:D:137:PRO:HD2	2.24	0.52
1:D:178:LEU:O	1:D:182:GLN:HB3	2.10	0.52
1:B:72:LEU:HD23	1:B:76:GLN:CG	2.40	0.52
1:C:282:ILE:HD12	1:D:282:ILE:HD12	1.92	0.52
1:B:66:ASN:OD1	2:F:508:VAL:HG13	2.10	0.52
1:A:284:ASN:ND2	1:B:284:ASN:ND2	2.55	0.52
1:A:212:MET:O	1:A:215:LEU:N	2.40	0.52
1:B:201:ILE:HG23	1:B:201:ILE:O	2.10	0.52
1:A:167:THR:HG23	1:A:168:ARG:N	2.25	0.52
1:A:180:GLU:O	1:A:181:CYS:HB2	2.09	0.52
1:D:101:LEU:HD23	1:D:101:LEU:O	2.10	0.51
1:C:133:GLY:O	1:C:135:ASP:N	2.36	0.51
1:A:219:VAL:HG22	1:A:219:VAL:O	2.10	0.51
1:D:201:ILE:O	1:D:201:ILE:HG23	2.10	0.51
1:B:204:LYS:HZ3	1:B:205:HIS:CE1	2.28	0.51
1:C:189:GLU:HB3	1:C:194:ILE:HG23	1.92	0.51
1:B:251:ILE:N	1:B:251:ILE:HD12	2.25	0.51
1:B:93:GLN:HG3	1:C:93:GLN:HE21	1.76	0.51
1:A:230:ASP:C	1:A:232:LEU:H	2.13	0.51
1:B:76:GLN:N	1:B:77:PRO:CD	2.73	0.51
1:A:133:GLY:O	1:A:135:ASP:N	2.35	0.51
1:C:101:LEU:HG	1:C:105:GLN:NE2	2.25	0.51
1:D:76:GLN:N	1:D:77:PRO:CD	2.73	0.51
1:B:29:GLU:C	1:B:31:LYS:H	2.14	0.51
1:D:72:LEU:HD23	1:D:76:GLN:CG	2.41	0.51
1:A:207:LYS:HB2	1:A:207:LYS:NZ	2.26	0.51
1:A:209:MET:O	1:A:212:MET:HB3	2.11	0.51
1:C:180:GLU:O	1:C:181:CYS:HB2	2.10	0.51
1:A:37:LEU:HD23	1:C:285:LEU:HD11	1.94	0.50
1:C:260:GLU:HG3	1:C:261:HIS:N	2.27	0.50
1:C:44:PRO:HG2	1:C:45:GLU:OE1	2.11	0.50
1:D:235:LYS:N	1:D:235:LYS:HE2	2.26	0.50
1:A:24:SER:N	1:A:25:PRO:CD	2.74	0.50
1:C:173:GLU:O	1:C:177:GLU:HG2	2.11	0.50
1:D:24:SER:HB2	1:D:25:PRO:HD3	1.92	0.50
1:D:76:GLN:O	1:D:77:PRO:C	2.49	0.50
1:D:273:ARG:HG2	1:D:273:ARG:HH11	1.76	0.50
1:D:251:ILE:N	1:D:251:ILE:HD12	2.26	0.50
1:B:56:ARG:HG2	1:B:56:ARG:HH11	1.77	0.50
1:B:37:LEU:HD23	1:B:42:LYS:HE2	1.94	0.50
1:A:173:GLU:O	1:A:177:GLU:HG2	2.11	0.50
1:D:26:SER:HB3	1:D:29:GLU:CD	2.32	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:30:LEU:HD13	1:A:52:ARG:HB2	1.94	0.50
1:A:42:LYS:HE3	1:D:249:SER:O	2.11	0.50
1:A:162:LEU:HD21	1:B:212:MET:HA	1.93	0.50
1:C:209:MET:O	1:C:212:MET:HB3	2.12	0.50
1:B:76:GLN:O	1:B:77:PRO:C	2.50	0.49
1:D:29:GLU:C	1:D:31:LYS:N	2.65	0.49
1:D:131:ASN:N	1:D:131:ASN:HD22	1.97	0.49
1:C:253:TYR:CD1	1:C:258:ILE:HG21	2.47	0.49
1:C:301:VAL:O	1:C:303:ASP:N	2.44	0.49
1:C:230:ASP:C	1:C:232:LEU:H	2.15	0.49
1:B:101:LEU:HD23	1:B:101:LEU:O	2.12	0.49
1:C:207:LYS:NZ	1:C:207:LYS:HB2	2.27	0.49
1:D:37:LEU:HD13	1:D:45:GLU:HB3	1.93	0.49
1:C:245:CYS:HA	1:C:284:ASN:H	1.77	0.49
1:A:260:GLU:HG3	1:A:261:HIS:N	2.28	0.49
1:A:215:LEU:CD2	1:B:219:VAL:HG21	2.42	0.49
1:D:145:LYS:HD3	1:D:304:TYR:HB3	1.94	0.49
1:B:235:LYS:N	1:B:235:LYS:HE2	2.28	0.49
1:D:37:LEU:HD23	1:D:42:LYS:HE2	1.95	0.49
1:B:129:ARG:HG2	1:B:129:ARG:HH11	1.78	0.49
1:A:301:VAL:O	1:A:303:ASP:N	2.45	0.49
1:A:259:GLU:HA	1:A:259:GLU:OE2	2.13	0.49
1:C:24:SER:N	1:C:25:PRO:CD	2.76	0.49
1:B:255:ARG:HA	1:B:258:ILE:HD11	1.95	0.49
1:D:56:ARG:HH11	1:D:56:ARG:HG2	1.78	0.49
1:C:235:LYS:HB2	1:C:253:TYR:CD2	2.48	0.49
1:B:164:SER:O	1:B:168:ARG:HG3	2.13	0.49
1:A:212:MET:HE2	1:B:166:LEU:HD21	1.95	0.48
1:D:164:SER:O	1:D:168:ARG:HG3	2.13	0.48
1:A:253:TYR:CD1	1:A:258:ILE:HG21	2.48	0.48
1:B:37:LEU:HD13	1:B:45:GLU:HB3	1.94	0.48
1:A:294:PHE:CD1	1:A:298:ASN:ND2	2.81	0.48
1:A:36:ARG:NH1	3:A:1307:SO4:S	2.86	0.48
1:D:76:GLN:HB2	1:D:80:ALA:HB2	1.95	0.48
1:B:76:GLN:HB2	1:B:80:ALA:HB2	1.95	0.48
1:A:182:GLN:CG	1:A:198:GLN:HE21	2.27	0.48
1:D:26:SER:HA	1:D:56:ARG:NH1	2.22	0.48
1:C:50:TYR:OH	2:G:508:VAL:HG11	2.14	0.48
1:C:286:ALA:C	1:D:287:MET:HE1	2.33	0.48
1:A:245:CYS:HA	1:A:284:ASN:H	1.77	0.48
1:D:286:ALA:O	1:D:290:VAL:HG23	2.13	0.48
1:D:255:ARG:HA	1:D:258:ILE:HD11	1.95	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:131:ASN:ND2	1:B:131:ASN:H	2.04	0.48
1:B:182:GLN:CG	1:B:198:GLN:HE22	2.26	0.48
1:D:164:SER:HA	1:D:167:THR:HG22	1.96	0.48
1:C:30:LEU:HD13	1:C:52:ARG:HB2	1.96	0.48
1:D:148:TRP:CE2	1:D:300:TRP:HA	2.49	0.48
1:D:242:ARG:HH11	1:D:242:ARG:HG2	1.79	0.48
1:B:145:LYS:HD3	1:B:304:TYR:HB3	1.94	0.48
1:B:286:ALA:O	1:B:290:VAL:HG23	2.13	0.48
1:D:30:LEU:HB2	1:D:53:ALA:HB2	1.96	0.48
1:B:241:MET:HE2	1:B:254:ASP:N	2.28	0.48
1:A:235:LYS:HB2	1:A:253:TYR:CD2	2.49	0.48
1:D:94:SER:HB3	1:D:97:ALA:CB	2.43	0.48
1:B:94:SER:HB3	1:B:97:ALA:CB	2.43	0.48
1:A:101:LEU:HG	1:A:105:GLN:NE2	2.29	0.48
1:C:248:PRO:HA	1:D:282:ILE:CD1	2.40	0.48
1:B:263:GLN:O	1:B:264:ARG:HB2	2.13	0.48
1:B:201:ILE:HD12	1:B:204:LYS:HZ2	1.79	0.48
1:B:242:ARG:HH11	1:B:242:ARG:HG2	1.79	0.48
1:C:258:ILE:HD13	1:C:281:LEU:HD11	1.96	0.47
1:D:132:PHE:CE1	2:H:505:MET:HG2	2.49	0.47
1:B:278:GLN:HG3	1:B:279:GLU:N	2.29	0.47
1:D:129:ARG:HG2	1:D:129:ARG:HH11	1.79	0.47
1:C:259:GLU:OE2	1:C:259:GLU:HA	2.14	0.47
1:A:84:CYS:C	1:A:86:ARG:H	2.17	0.47
1:B:29:GLU:C	1:B:31:LYS:N	2.68	0.47
1:C:38:PHE:CZ	1:C:69:LEU:HG	2.50	0.47
1:A:138:SER:HA	1:A:238:PHE:CD1	2.50	0.47
1:D:263:GLN:O	1:D:264:ARG:HB2	2.13	0.47
1:B:26:SER:HA	1:B:56:ARG:NH1	2.23	0.47
1:D:131:ASN:H	1:D:131:ASN:ND2	2.05	0.47
1:C:229:PRO:HB3	1:C:231:TYR:CE2	2.49	0.47
1:C:294:PHE:CD1	1:C:298:ASN:ND2	2.83	0.47
1:A:44:PRO:HG2	1:A:45:GLU:OE1	2.14	0.47
1:C:129:ARG:NH2	1:C:264:ARG:HH21	2.12	0.47
1:D:223:ARG:HG2	1:D:223:ARG:NH1	2.30	0.47
1:B:148:TRP:CE2	1:B:300:TRP:HA	2.50	0.47
1:D:133:GLY:O	1:D:135:ASP:N	2.47	0.47
1:D:223:ARG:HG2	1:D:223:ARG:HH11	1.80	0.47
1:D:278:GLN:HG3	1:D:279:GLU:N	2.29	0.47
1:A:239:GLU:HG2	1:A:240:LEU:N	2.30	0.47
1:C:84:CYS:C	1:C:86:ARG:H	2.17	0.47
1:C:176:ARG:HH12	1:D:173:GLU:HG2	1.80	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:138:SER:O	1:A:142:ILE:HG13	2.15	0.47
1:A:199:ALA:HA	1:A:202:GLU:HB2	1.97	0.47
1:A:43:TYR:N	1:A:44:PRO:HD2	2.30	0.46
1:C:182:GLN:CG	1:C:198:GLN:HE21	2.28	0.46
1:D:241:MET:HE2	1:D:254:ASP:N	2.30	0.46
1:C:71:TYR:CD2	1:C:79:GLN:HB2	2.51	0.46
1:C:138:SER:O	1:C:142:ILE:HG13	2.15	0.46
1:C:111:TYR:HB3	1:C:147:ARG:HG3	1.96	0.46
1:A:229:PRO:HB3	1:A:231:TYR:CE2	2.50	0.46
1:B:223:ARG:HG2	1:B:223:ARG:NH1	2.30	0.46
1:A:257:ASP:C	1:A:259:GLU:H	2.19	0.46
1:D:105:GLN:C	1:D:107:GLU:H	2.17	0.46
1:B:105:GLN:C	1:B:107:GLU:H	2.18	0.46
1:C:167:THR:CG2	1:C:168:ARG:N	2.78	0.46
1:A:38:PHE:CE1	2:E:508:VAL:HG23	2.50	0.46
1:D:154:ARG:HH11	1:D:154:ARG:HG3	1.80	0.46
1:C:172:ALA:O	1:C:176:ARG:HG3	2.16	0.46
1:C:81:LEU:O	1:C:81:LEU:HD13	2.16	0.46
1:D:76:GLN:OE1	1:D:76:GLN:N	2.45	0.46
1:C:66:ASN:OD1	2:G:508:VAL:HG13	2.15	0.46
1:B:164:SER:HA	1:B:167:THR:HG22	1.98	0.46
1:B:223:ARG:HH11	1:B:223:ARG:HG2	1.81	0.46
1:C:239:GLU:HG2	1:C:240:LEU:N	2.31	0.46
1:A:211:ASP:HA	1:A:214:GLU:OE1	2.16	0.46
1:C:257:ASP:C	1:C:259:GLU:H	2.19	0.46
1:B:273:ARG:NH1	1:B:273:ARG:HG2	2.29	0.46
1:A:26:SER:OG	1:A:29:GLU:HG3	2.15	0.46
1:A:38:PHE:CZ	1:A:69:LEU:HG	2.51	0.46
1:B:294:PHE:CE1	1:B:298:ASN:ND2	2.84	0.46
1:D:157:HIS:O	1:D:161:GLU:HB2	2.15	0.46
1:C:211:ASP:HA	1:C:214:GLU:OE1	2.16	0.46
1:D:169:LEU:O	1:D:173:GLU:HG3	2.15	0.45
1:D:73:LYS:O	1:D:74:MET:HB3	2.17	0.45
1:B:182:GLN:NE2	1:B:198:GLN:NE2	2.63	0.45
1:D:182:GLN:CG	1:D:198:GLN:HE22	2.29	0.45
1:A:47:ALA:HB2	1:A:70:CYS:HB3	1.98	0.45
1:A:129:ARG:NH2	1:A:264:ARG:HH21	2.14	0.45
1:D:38:PHE:O	1:D:41:ARG:N	2.50	0.45
1:D:254:ASP:O	1:D:256:LYS:N	2.48	0.45
1:A:167:THR:CG2	1:A:168:ARG:N	2.79	0.45
1:C:138:SER:HA	1:C:238:PHE:CD1	2.51	0.45
1:B:154:ARG:HH11	1:B:154:ARG:HG3	1.80	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:76:GLN:OE1	1:B:76:GLN:N	2.46	0.45
1:D:258:ILE:CD1	1:D:259:GLU:N	2.74	0.45
1:C:212:MET:HE2	1:D:166:LEU:HD21	1.99	0.45
1:B:73:LYS:O	1:B:74:MET:HB3	2.17	0.45
1:B:258:ILE:CD1	1:B:259:GLU:N	2.74	0.45
1:A:258:ILE:HD13	1:A:281:LEU:HD11	1.99	0.45
1:B:133:GLY:O	1:B:135:ASP:N	2.49	0.45
2:H:502:THR:HG22	2:H:503:SER:N	2.32	0.45
1:A:172:ALA:O	1:A:176:ARG:HG3	2.17	0.45
1:C:32:GLU:O	1:C:35:ASN:HB2	2.17	0.45
1:A:42:LYS:HD3	1:C:283:PRO:HG3	1.97	0.45
1:A:110:SER:HB3	1:A:113:GLU:OE1	2.16	0.45
1:B:254:ASP:O	1:B:256:LYS:N	2.49	0.45
1:B:46:ALA:HA	1:B:49:CYS:HB3	1.99	0.45
1:C:176:ARG:NH1	1:D:173:GLU:HG2	2.32	0.45
1:B:57:ASN:ND2	1:B:57:ASN:C	2.69	0.45
2:E:502:THR:HG22	2:E:503:SER:N	2.31	0.45
1:D:226:ARG:HD3	1:D:300:TRP:CZ3	2.52	0.45
1:C:199:ALA:HA	1:C:202:GLU:HB2	1.99	0.45
1:B:169:LEU:O	1:B:173:GLU:HG3	2.16	0.45
1:D:294:PHE:CE1	1:D:298:ASN:ND2	2.85	0.45
1:D:156:ILE:HD12	1:D:156:ILE:N	2.32	0.45
1:B:69:LEU:HD12	1:B:69:LEU:O	2.17	0.45
1:A:141:ARG:HG3	1:A:141:ARG:NH1	2.31	0.45
1:D:46:ALA:HA	1:D:49:CYS:HB3	1.99	0.45
1:D:182:GLN:O	1:D:184:ASN:N	2.50	0.45
1:A:85:ARG:HH11	1:A:85:ARG:CG	2.29	0.45
1:D:242:ARG:CG	1:D:242:ARG:HH11	2.30	0.45
1:A:81:LEU:O	1:A:81:LEU:HD13	2.17	0.45
1:B:132:PHE:O	1:B:135:ASP:HB2	2.17	0.44
1:B:26:SER:HB3	1:B:29:GLU:CG	2.48	0.44
1:B:253:TYR:HB2	1:B:258:ILE:HG12	1.99	0.44
1:D:253:TYR:HB2	1:D:258:ILE:HG12	1.99	0.44
1:B:69:LEU:HD23	2:F:508:VAL:HG22	1.97	0.44
1:B:115:ILE:HD13	1:B:144:LYS:HD3	1.99	0.44
1:A:32:GLU:O	1:A:35:ASN:HB2	2.17	0.44
1:C:110:SER:HB3	1:C:113:GLU:OE1	2.17	0.44
1:B:156:ILE:HD12	1:B:156:ILE:N	2.32	0.44
1:C:141:ARG:NH1	1:C:141:ARG:HG3	2.31	0.44
1:A:169:LEU:HD22	1:B:173:GLU:CD	2.37	0.44
1:D:204:LYS:NZ	1:D:205:HIS:HE1	2.15	0.44
1:D:132:PHE:O	1:D:135:ASP:HB2	2.17	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:134:ASP:O	1:D:138:SER:CB	2.65	0.44
1:B:134:ASP:O	1:B:138:SER:CB	2.66	0.44
2:F:502:THR:HG22	2:F:503:SER:N	2.33	0.44
1:D:57:ASN:ND2	1:D:57:ASN:C	2.70	0.44
1:A:30:LEU:HB2	1:A:53:ALA:HB2	1.99	0.44
1:D:115:ILE:HD13	1:D:144:LYS:HD3	1.99	0.44
1:D:226:ARG:HD3	1:D:300:TRP:CE3	2.52	0.44
1:B:132:PHE:CE1	2:F:505:MET:HG2	2.52	0.44
1:B:131:ASN:CB	2:F:502:THR:HG23	2.48	0.44
1:B:242:ARG:HH11	1:B:242:ARG:CG	2.30	0.44
1:D:182:GLN:OE1	1:D:183:ARG:HB2	2.17	0.44
1:C:47:ALA:HB2	1:C:70:CYS:HB3	1.99	0.44
1:D:182:GLN:C	1:D:184:ASN:H	2.20	0.44
1:A:174:ARG:O	1:A:178:LEU:HD13	2.18	0.44
1:B:204:LYS:HZ3	1:B:205:HIS:HE1	1.64	0.44
1:C:105:GLN:OE1	1:C:113:GLU:HB3	2.17	0.44
1:C:85:ARG:CG	1:C:85:ARG:HH11	2.30	0.44
2:G:502:THR:HG22	2:G:503:SER:N	2.32	0.44
1:D:273:ARG:NH1	1:D:273:ARG:HG2	2.31	0.44
1:B:157:HIS:O	1:B:161:GLU:HB2	2.17	0.44
1:C:95:VAL:HG11	1:C:125:ALA:HB2	2.00	0.44
1:C:237:SER:O	1:C:238:PHE:HB2	2.18	0.43
1:A:287:MET:O	1:A:288:LYS:C	2.56	0.43
1:D:247:THR:HB	1:D:248:PRO:CD	2.48	0.43
1:C:163:HIS:O	1:C:165:TYR:N	2.52	0.43
1:A:71:TYR:CD2	1:A:79:GLN:HB2	2.53	0.43
1:B:204:LYS:NZ	1:B:205:HIS:HE1	2.16	0.43
1:B:226:ARG:HD3	1:B:300:TRP:CE3	2.53	0.43
1:B:131:ASN:CG	2:F:502:THR:HG23	2.39	0.43
1:C:37:LEU:HA	1:C:37:LEU:HD23	1.82	0.43
1:D:76:GLN:CD	1:D:76:GLN:N	2.70	0.43
1:B:182:GLN:O	1:B:184:ASN:N	2.52	0.43
1:B:226:ARG:HD3	1:B:300:TRP:CZ3	2.53	0.43
1:C:26:SER:OG	1:C:29:GLU:HG3	2.17	0.43
1:C:195:ARG:HH11	1:C:195:ARG:HG3	1.83	0.43
1:B:239:GLU:O	1:B:240:LEU:C	2.57	0.43
2:H:507:GLU:HG3	2:H:507:GLU:O	2.19	0.43
1:B:182:GLN:C	1:B:184:ASN:H	2.21	0.43
1:C:215:LEU:CD2	1:D:219:VAL:HG21	2.48	0.43
1:C:174:ARG:O	1:C:178:LEU:HD13	2.19	0.43
1:C:30:LEU:HB2	1:C:53:ALA:HB2	2.01	0.43
1:B:77:PRO:HB2	1:B:78:GLU:H	1.63	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:F:507:GLU:O	2:F:507:GLU:HG3	2.19	0.43
1:B:243:GLU:HB3	1:B:255:ARG:HB2	1.99	0.43
1:D:254:ASP:C	1:D:256:LYS:H	2.22	0.43
1:B:182:GLN:HG3	1:B:198:GLN:CD	2.38	0.43
1:A:111:TYR:HB3	1:A:147:ARG:HG3	2.00	0.43
1:D:77:PRO:HB2	1:D:78:GLU:H	1.63	0.43
1:C:284:ASN:ND2	1:D:284:ASN:ND2	2.62	0.43
1:B:30:LEU:HB2	1:B:53:ALA:HB2	2.01	0.43
1:B:134:ASP:O	1:B:138:SER:HB2	2.19	0.43
1:B:182:GLN:OE1	1:B:183:ARG:HB2	2.19	0.43
1:C:235:LYS:HD3	1:C:235:LYS:HA	1.85	0.43
1:A:81:LEU:HD22	1:A:101:LEU:HD12	2.00	0.43
1:B:247:THR:HB	1:B:248:PRO:CD	2.49	0.43
1:D:182:GLN:HG3	1:D:198:GLN:CD	2.38	0.43
1:A:235:LYS:HA	1:A:235:LYS:HD3	1.82	0.43
1:C:96:LYS:HA	2:G:505:MET:HE1	2.01	0.43
1:C:81:LEU:HD22	1:C:101:LEU:HD12	2.01	0.43
1:C:258:ILE:HD13	1:C:281:LEU:CD1	2.49	0.43
1:C:287:MET:O	1:C:288:LYS:C	2.57	0.43
1:B:228:ILE:HD12	1:B:294:PHE:HE2	1.84	0.43
1:D:170:ILE:HG22	1:D:171:ALA:N	2.33	0.43
1:B:254:ASP:C	1:B:256:LYS:H	2.22	0.42
1:D:294:PHE:O	1:D:298:ASN:ND2	2.52	0.42
1:C:43:TYR:N	1:C:44:PRO:HD2	2.34	0.42
1:B:61:ALA:O	1:B:62:VAL:C	2.58	0.42
1:B:76:GLN:CD	1:B:76:GLN:N	2.70	0.42
1:A:50:TYR:OH	2:E:508:VAL:HG11	2.19	0.42
1:D:239:GLU:O	1:D:240:LEU:C	2.57	0.42
1:C:178:LEU:N	1:C:178:LEU:HD12	2.35	0.42
1:A:195:ARG:HG3	1:A:195:ARG:HH11	1.84	0.42
1:B:215:LEU:O	1:B:218:GLN:HB2	2.20	0.42
1:B:241:MET:CE	1:B:254:ASP:N	2.82	0.42
1:A:230:ASP:C	1:A:232:LEU:N	2.73	0.42
1:B:203:ALA:HA	1:B:206:ASP:OD2	2.19	0.42
1:C:284:ASN:OD1	1:C:287:MET:HB2	2.18	0.42
1:B:145:LYS:NZ	1:B:304:TYR:HB3	2.35	0.42
1:C:176:ARG:HA	1:C:180:GLU:HB2	2.01	0.42
1:C:57:ASN:C	1:C:57:ASN:HD22	2.21	0.42
1:B:145:LYS:CD	1:B:304:TYR:HB3	2.50	0.42
1:A:178:LEU:HD12	1:A:178:LEU:N	2.35	0.42
1:D:247:THR:HB	1:D:248:PRO:HD2	2.00	0.42
1:D:182:GLN:NE2	1:D:198:GLN:NE2	2.67	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:113:GLU:O	1:D:116:ALA:HB3	2.19	0.42
1:B:69:LEU:HD23	2:F:508:VAL:HG23	2.01	0.42
1:A:154:ARG:O	1:A:154:ARG:HG3	2.19	0.42
1:A:37:LEU:HD23	1:A:37:LEU:HA	1.82	0.42
1:B:170:ILE:HG22	1:B:171:ALA:N	2.33	0.42
1:D:243:GLU:HB3	1:D:255:ARG:HB2	2.00	0.42
1:A:176:ARG:HA	1:A:180:GLU:HB2	2.01	0.42
1:C:154:ARG:O	1:C:154:ARG:HG3	2.20	0.42
1:D:228:ILE:HD12	1:D:294:PHE:HE2	1.85	0.42
1:C:163:HIS:O	1:C:164:SER:C	2.57	0.42
1:D:85:ARG:CG	1:D:85:ARG:HH11	2.32	0.42
1:A:237:SER:O	1:A:238:PHE:HB2	2.20	0.42
1:A:31:LYS:HE2	1:A:31:LYS:HB3	1.86	0.42
1:D:178:LEU:C	1:D:180:GLU:H	2.23	0.41
1:C:169:LEU:HD22	1:D:173:GLU:CD	2.41	0.41
1:C:113:GLU:O	1:C:116:ALA:HB3	2.19	0.41
1:D:107:GLU:HA	1:D:107:GLU:OE1	2.20	0.41
1:D:38:PHE:O	1:D:40:GLY:N	2.53	0.41
1:B:254:ASP:O	1:B:258:ILE:HD11	2.20	0.41
1:C:38:PHE:CE1	2:G:508:VAL:HG23	2.54	0.41
1:B:85:ARG:CG	1:B:85:ARG:HH11	2.32	0.41
1:D:148:TRP:O	1:D:150:SER:N	2.53	0.41
1:A:105:GLN:OE1	1:A:113:GLU:HB3	2.19	0.41
1:A:31:LYS:HD3	1:A:63:TYR:HE1	1.85	0.41
1:A:258:ILE:HD13	1:A:281:LEU:CD1	2.50	0.41
1:D:148:TRP:CZ2	1:D:300:TRP:HB3	2.55	0.41
1:A:113:GLU:O	1:A:116:ALA:HB3	2.20	0.41
1:A:95:VAL:HG11	1:A:125:ALA:HB2	2.02	0.41
1:D:61:ALA:O	1:D:62:VAL:C	2.59	0.41
1:A:163:HIS:O	1:A:164:SER:C	2.58	0.41
1:A:163:HIS:O	1:A:165:TYR:N	2.54	0.41
1:D:26:SER:HB3	1:D:29:GLU:CG	2.51	0.41
1:B:178:LEU:C	1:B:180:GLU:H	2.24	0.41
1:C:268:PHE:HA	1:C:274:SER:O	2.21	0.41
1:C:235:LYS:HB2	1:C:253:TYR:CE2	2.56	0.41
1:B:166:LEU:O	1:B:169:LEU:N	2.53	0.41
1:A:248:PRO:HA	1:B:282:ILE:CD1	2.46	0.41
1:A:268:PHE:HA	1:A:274:SER:O	2.21	0.41
1:D:134:ASP:O	1:D:138:SER:HB2	2.21	0.41
1:D:254:ASP:O	1:D:258:ILE:HD11	2.21	0.41
1:C:248:PRO:HD2	1:C:276:LEU:HD13	2.03	0.41
1:D:145:LYS:CD	1:D:304:TYR:HB3	2.50	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:107:GLU:HA	1:B:107:GLU:OE1	2.21	0.41
1:D:288:LYS:HE3	1:D:288:LYS:HB2	1.87	0.41
1:D:255:ARG:HG3	1:D:255:ARG:O	2.21	0.41
1:A:248:PRO:HD2	1:A:276:LEU:HD13	2.03	0.41
1:A:134:ASP:O	1:A:135:ASP:C	2.57	0.41
1:B:140:LEU:O	1:B:141:ARG:C	2.58	0.41
1:A:96:LYS:HA	2:E:505:MET:HE1	2.02	0.41
2:E:507:GLU:O	2:E:507:GLU:HG3	2.21	0.41
1:A:37:LEU:HD13	1:A:45:GLU:HB3	2.03	0.41
1:C:230:ASP:C	1:C:232:LEU:N	2.75	0.41
1:B:247:THR:HB	1:B:248:PRO:HD2	2.02	0.41
1:D:215:LEU:O	1:D:218:GLN:HB2	2.21	0.41
1:D:131:ASN:O	1:D:131:ASN:ND2	2.54	0.41
1:C:251:ILE:N	1:C:251:ILE:CD1	2.78	0.41
1:A:156:ILE:C	1:A:158:GLN:H	2.25	0.41
1:D:241:MET:CE	1:D:254:ASP:N	2.84	0.40
1:B:38:PHE:O	1:B:40:GLY:N	2.54	0.40
1:B:37:LEU:HD12	1:B:46:ALA:HA	2.03	0.40
1:C:156:ILE:C	1:C:158:GLN:H	2.25	0.40
1:D:203:ALA:HA	1:D:206:ASP:OD2	2.21	0.40
1:C:71:TYR:CG	1:C:79:GLN:HB2	2.57	0.40
1:D:85:ARG:NH1	1:D:85:ARG:HG3	2.35	0.40
1:A:136:ILE:N	1:A:137:PRO:HD2	2.36	0.40
1:A:194:ILE:O	1:A:198:GLN:HB2	2.22	0.40
1:C:302:GLU:OE2	1:C:302:GLU:HA	2.22	0.40
1:A:141:ARG:HG3	1:A:141:ARG:HH11	1.86	0.40
1:B:241:MET:SD	1:B:252:THR:HG22	2.61	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	279/281 (99%)	223 (80%)	45 (16%)	11 (4%)	5 37

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	279/281 (99%)	214 (77%)	51 (18%)	14 (5%)	3	28
1	C	279/281 (99%)	223 (80%)	46 (16%)	10 (4%)	5	40
1	D	279/281 (99%)	216 (77%)	48 (17%)	15 (5%)	3	26
2	E	7/9 (78%)	3 (43%)	2 (29%)	2 (29%)	0	0
2	F	7/9 (78%)	3 (43%)	2 (29%)	2 (29%)	0	0
2	G	7/9 (78%)	3 (43%)	2 (29%)	2 (29%)	0	0
2	H	7/9 (78%)	3 (43%)	2 (29%)	2 (29%)	0	0
All	All	1144/1160 (99%)	888 (78%)	198 (17%)	58 (5%)	3	28

All (58) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	159	GLU
1	A	181	CYS
1	A	300	TRP
1	A	301	VAL
1	B	38	PHE
1	B	77	PRO
1	B	134	ASP
1	B	182	GLN
1	B	264	ARG
1	C	159	GLU
1	C	181	CYS
1	C	300	TRP
1	C	301	VAL
1	D	38	PHE
1	D	77	PRO
1	D	134	ASP
1	D	182	GLN
1	D	264	ARG
2	E	502	THR
2	F	502	THR
2	G	502	THR
2	H	502	THR
1	B	74	MET
1	B	75	GLN
1	B	180	GLU
1	D	74	MET
1	D	75	GLN
1	D	180	GLU

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Mol	Chain	Res	Type
2	E	503	SER
2	F	503	SER
2	G	503	SER
2	H	503	SER
1	A	133	GLY
1	A	158	GLN
1	A	160	SER
1	B	30	LEU
1	B	227	ASP
1	B	255	ARG
1	C	158	GLN
1	C	160	SER
1	D	30	LEU
1	D	183	ARG
1	D	227	ASP
1	D	255	ARG
1	B	183	ARG
1	C	133	GLY
1	A	190	ASP
1	B	178	LEU
1	C	164	SER
1	C	190	ASP
1	A	111	TYR
1	A	164	SER
1	D	178	LEU
1	A	270	PRO
1	C	270	PRO
1	B	76	GLN
1	D	76	GLN
1	D	39	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	244/244 (100%)	224 (92%)	20 (8%)	17	57
1	B	244/244 (100%)	220 (90%)	24 (10%)	12	45

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	244/244 (100%)	223 (91%)	21 (9%)	15	54
1	D	244/244 (100%)	220 (90%)	24 (10%)	12	45
2	E	9/9 (100%)	8 (89%)	1 (11%)	9	37
2	F	9/9 (100%)	9 (100%)	0	100	100
2	G	9/9 (100%)	9 (100%)	0	100	100
2	H	9/9 (100%)	8 (89%)	1 (11%)	9	37
All	All	1012/1012 (100%)	921 (91%)	91 (9%)	14	50

All (91) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	CYS
1	A	55	THR
1	A	57	ASN
1	A	103	GLN
1	A	135	ASP
1	A	141	ARG
1	A	149	ASN
1	A	151	ILE
1	A	195	ARG
1	A	198	GLN
1	A	204	LYS
1	A	208	TYR
1	A	227	ASP
1	A	240	LEU
1	A	242	ARG
1	A	257	ASP
1	A	273	ARG
1	A	280	GLN
1	A	288	LYS
1	A	289	GLU
1	B	24	SER
1	B	32	GLU
1	B	41	ARG
1	B	55	THR
1	B	57	ASN
1	B	75	GLN
1	B	76	GLN
1	B	104	CYS
1	B	106	LEU

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Mol	Chain	Res	Type
1	B	131	ASN
1	B	135	ASP
1	B	146	LYS
1	B	154	ARG
1	B	161	GLU
1	B	166	LEU
1	B	175	GLU
1	B	195	ARG
1	B	198	GLN
1	B	213	ASP
1	B	223	ARG
1	B	258	ILE
1	B	273	ARG
1	B	278	GLN
1	B	292	ASP
1	C	49	CYS
1	C	55	THR
1	C	57	ASN
1	C	83	ASP
1	C	103	GLN
1	C	135	ASP
1	C	141	ARG
1	C	149	ASN
1	C	151	ILE
1	C	195	ARG
1	C	198	GLN
1	C	204	LYS
1	C	208	TYR
1	C	227	ASP
1	C	240	LEU
1	C	242	ARG
1	C	257	ASP
1	C	273	ARG
1	C	280	GLN
1	C	288	LYS
1	C	289	GLU
1	D	24	SER
1	D	32	GLU
1	D	41	ARG
1	D	55	THR
1	D	57	ASN
1	D	75	GLN

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Mol	Chain	Res	Type
1	D	76	GLN
1	D	104	CYS
1	D	106	LEU
1	D	131	ASN
1	D	135	ASP
1	D	146	LYS
1	D	154	ARG
1	D	161	GLU
1	D	166	LEU
1	D	175	GLU
1	D	195	ARG
1	D	198	GLN
1	D	213	ASP
1	D	223	ARG
1	D	258	ILE
1	D	273	ARG
1	D	278	GLN
1	D	292	ASP
2	E	509	ASP
2	H	509	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
1	A	66	ASN
1	A	79	GLN
1	A	117	ASN
1	A	131	ASN
1	A	182	GLN
1	A	198	GLN
1	A	205	HIS
1	A	218	GLN
1	A	280	GLN
1	A	284	ASN
1	A	298	ASN
1	B	28	GLN
1	B	57	ASN
1	B	93	GLN
1	B	131	ASN
1	B	184	ASN
1	B	198	GLN

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Mol	Chain	Res	Type
1	B	205	HIS
1	B	267	HIS
1	B	278	GLN
1	B	280	GLN
1	C	66	ASN
1	C	79	GLN
1	C	93	GLN
1	C	117	ASN
1	C	131	ASN
1	C	182	GLN
1	C	198	GLN
1	C	205	HIS
1	C	218	GLN
1	C	280	GLN
1	C	284	ASN
1	C	298	ASN
1	D	57	ASN
1	D	131	ASN
1	D	184	ASN
1	D	198	GLN
1	D	205	HIS
1	D	267	HIS
1	D	278	GLN
1	D	280	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 11 ligands modelled in this entry, 4 are monoatomic - leaving 7 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	SO4	A	1305	-	4,4,4	0.25	0	6,6,6	0.07	0
3	SO4	A	1306	-	4,4,4	0.41	0	6,6,6	0.08	0
3	SO4	A	1307	-	4,4,4	0.79	0	6,6,6	0.13	0
3	SO4	B	1305	-	4,4,4	0.32	0	6,6,6	0.12	0
3	SO4	B	1306	-	4,4,4	0.35	0	6,6,6	0.16	0
3	SO4	C	1305	-	4,4,4	0.25	0	6,6,6	0.15	0
3	SO4	D	1305	-	4,4,4	0.39	0	6,6,6	0.16	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	SO4	A	1305	-	-	0/0/0/0	0/0/0/0
3	SO4	A	1306	-	-	0/0/0/0	0/0/0/0
3	SO4	A	1307	-	-	0/0/0/0	0/0/0/0
3	SO4	B	1305	-	-	0/0/0/0	0/0/0/0
3	SO4	B	1306	-	-	0/0/0/0	0/0/0/0
3	SO4	C	1305	-	-	0/0/0/0	0/0/0/0
3	SO4	D	1305	-	-	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.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95th 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 ²)	Q<0.9
1	A	281/281 (100%)	0.22	2 (0%) 84 42	36, 74, 120, 136	0
1	B	281/281 (100%)	0.17	1 (0%) 90 57	36, 78, 117, 136	0
1	C	281/281 (100%)	0.19	2 (0%) 84 42	36, 74, 120, 135	0
1	D	281/281 (100%)	0.21	3 (1%) 77 30	36, 78, 117, 136	0
2	E	9/9 (100%)	0.27	0 100 100	95, 103, 114, 120	0
2	F	9/9 (100%)	0.62	0 100 100	95, 103, 115, 120	0
2	G	9/9 (100%)	0.27	0 100 100	96, 102, 115, 120	0
2	H	9/9 (100%)	0.63	0 100 100	95, 103, 116, 121	0
All	All	1160/1160 (100%)	0.21	8 (0%) 79 42	36, 78, 121, 136	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	182	GLN	3.2
1	C	182	GLN	2.8
1	D	189	GLU	2.8
1	D	188	HIS	2.7
1	A	189	GLU	2.5
1	B	189	GLU	2.3
1	C	180	GLU	2.3
1	D	186	GLU	2.2

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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	SO4	A	1307	5/5	0.60	5.66	108,109,111,112	0
3	SO4	A	1305	5/5	0.27	-0.21	89,90,90,91	0
3	SO4	B	1305	5/5	0.27	-0.67	82,83,85,86	0
3	SO4	B	1306	5/5	0.24	-0.90	89,89,91,91	0
3	SO4	A	1306	5/5	0.21	-1.00	78,79,80,80	0
3	SO4	D	1305	5/5	0.21	-1.41	105,106,107,107	0
3	SO4	C	1305	5/5	0.20	-1.47	88,89,90,91	0
4	NI	B	1307	1/1	0.52	-	200,200,200,200	0
4	NI	B	1308	1/1	0.46	-	200,200,200,200	0
4	NI	B	1309	1/1	0.42	-	200,200,200,200	0
4	NI	B	1310	1/1	0.29	-	200,200,200,200	0

6.5 Other polymers ⓘ

There are no such residues in this entry.