



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:17 AM UTC

PDB ID : 3A1Z / pdb_00003a1z
Title : Crystal structure of juvenile hormone binding protein from silkworm
Authors : Suzuki, R.; Fujimoto, Z.; Shiotsuki, T.; Momma, M.; Tase, A.; Yamazaki, T.
Deposited on : 2009-04-27
Resolution : 2.59 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

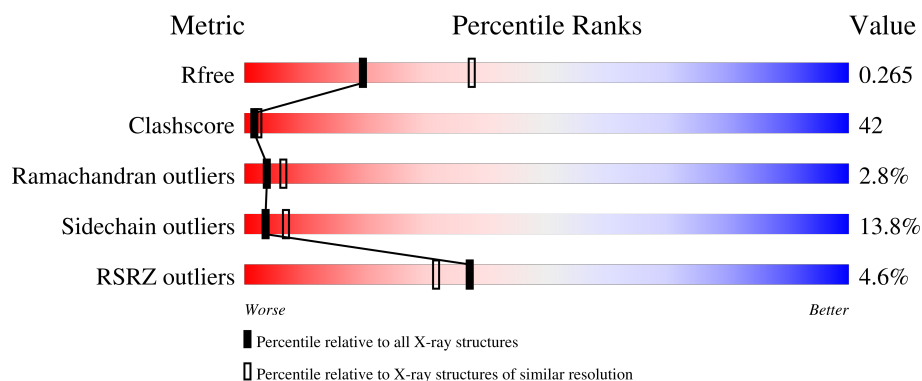
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	227	0% 45% 37% 10% 8%
1	B	227	7% 36% 44% 12% 7%
1	C	227	2% 51% 35% 7% 7%
1	D	227	6% 45% 40% 7% 8%

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	ZN	A	228	-	-	X	-
2	ZN	A	230	-	-	-	X
2	ZN	C	231	-	-	-	X
3	MPD	A	231	-	-	X	-
3	MPD	A	232	-	-	X	-
3	MPD	B	230	-	-	X	-
3	MPD	B	231	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6672 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 Hemolymph juvenile hormone binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	0	0	0
			1599	1014	262	316	7			
1	B	212	Total	C	N	O	S	0	0	0
			1630	1032	267	323	8			
1	C	212	Total	C	N	O	S	0	0	0
			1631	1034	268	322	7			
1	D	209	Total	C	N	O	S	0	0	0
			1607	1019	263	317	8			

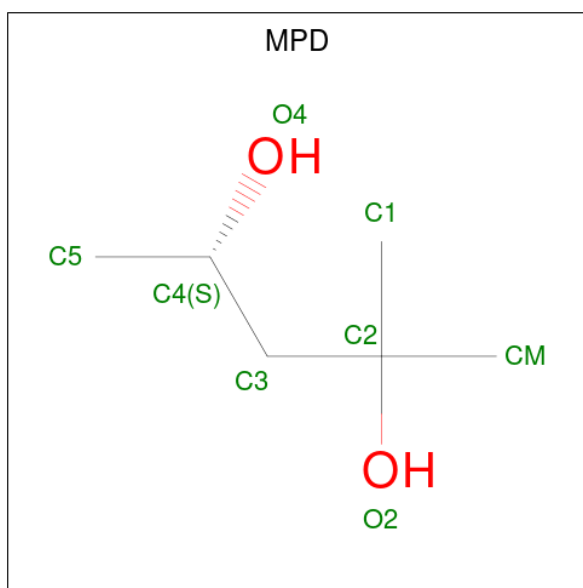
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP Q9U556
A	-1	SER	-	expression tag	UNP Q9U556
B	-2	GLY	-	expression tag	UNP Q9U556
B	-1	SER	-	expression tag	UNP Q9U556
C	-2	GLY	-	expression tag	UNP Q9U556
C	-1	SER	-	expression tag	UNP Q9U556
D	-2	GLY	-	expression tag	UNP Q9U556
D	-1	SER	-	expression tag	UNP Q9U556

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	5	Total	Zn	0	0
			5	5		
2	B	4	Total	Zn	0	0
			4	4		
2	C	6	Total	Zn	0	0
			6	6		
2	D	5	Total	Zn	0	0
			5	5		

- Molecule 3 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			8	6	2		
3	A	1	Total	C	O	0	0
			8	6	2		
3	B	1	Total	C	O	0	0
			8	6	2		
3	B	1	Total	C	O	0	0
			8	6	2		
3	C	1	Total	C	O	0	0
			8	6	2		
3	C	1	Total	C	O	0	0
			8	6	2		
3	D	1	Total	C	O	0	0
			8	6	2		
3	D	1	Total	C	O	0	0
			8	6	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	33	Total	O	0	0
			33	33		
4	B	24	Total	O	0	0
			24	24		
4	C	42	Total	O	0	0
			42	42		

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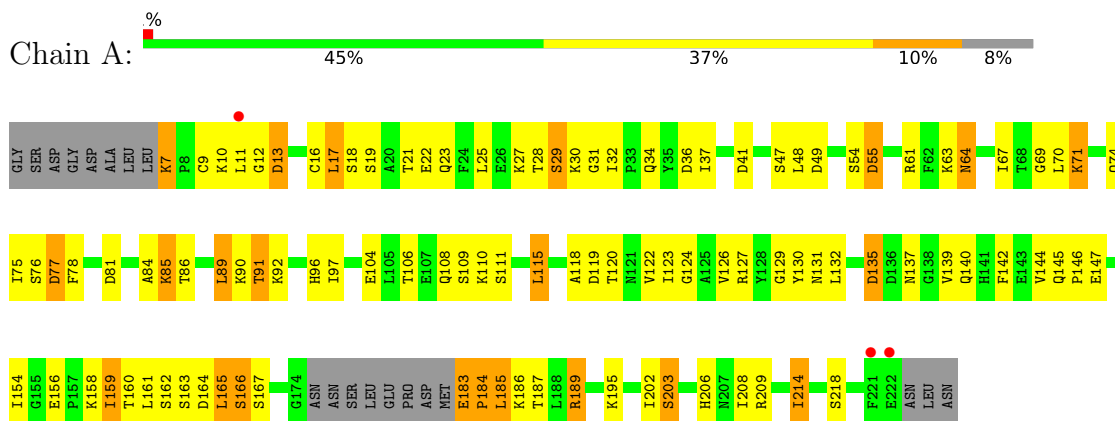
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	22	Total	O	0	0
			22	22		

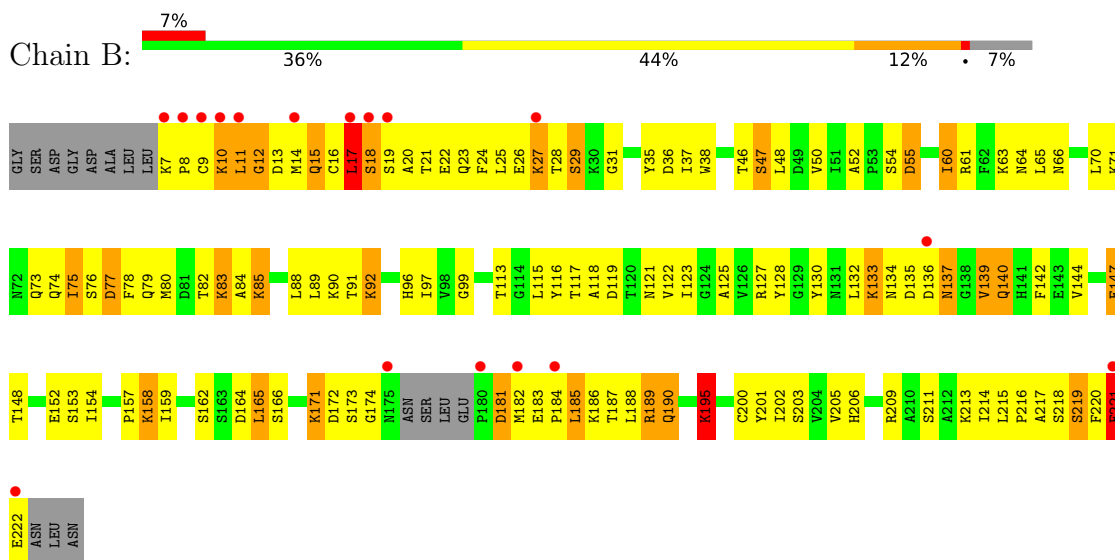
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

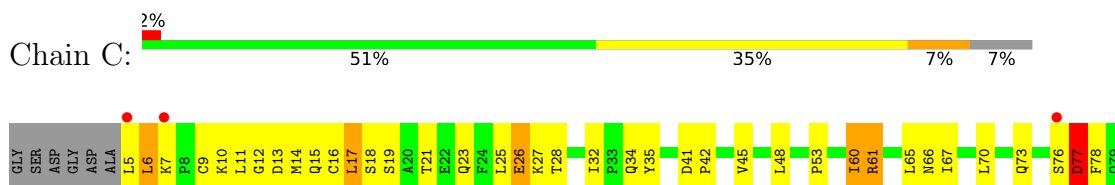
- Molecule 1: Hemolymph juvenile hormone binding protein

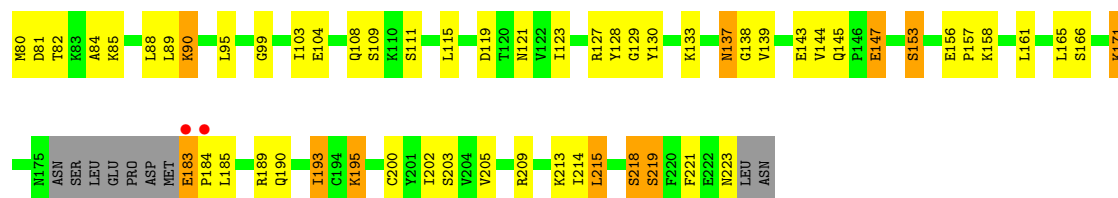


- Molecule 1: Hemolymph juvenile hormone binding protein

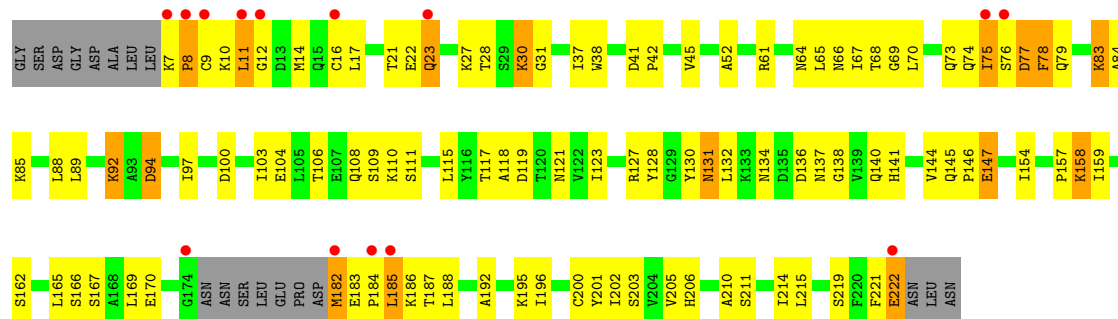


- Molecule 1: Hemolymph juvenile hormone binding protein





• Molecule 1: Hemolymph juvenile hormone binding protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	54.86Å 114.66Å 192.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.49 – 2.59 49.49 – 2.59	Depositor EDS
% Data completeness (in resolution range)	99.4 (49.49-2.59) 99.4 (49.49-2.59)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.89 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.5.0066	Depositor
R, R_{free}	0.222 , 0.290 (Not available) , 0.265	Depositor DCC
R_{free} test set	1933 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	51.9	Xtriage
Anisotropy	0.143	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6672	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 21.23 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 7.3350e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical 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: MPD, ZN

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	0.98	0/1624	1.17	8/2199 (0.4%)
1	B	0.87	0/1656	1.16	5/2242 (0.2%)
1	C	0.98	0/1656	1.18	3/2243 (0.1%)
1	D	0.92	0/1632	1.15	1/2209 (0.0%)
All	All	0.94	0/6568	1.16	17/8893 (0.2%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	202	ILE	N-CA-C	-8.75	101.69	110.62
1	A	202	ILE	N-CA-C	-7.70	102.77	110.62
1	C	26	GLU	N-CA-C	-6.59	103.79	110.97
1	B	27	LYS	N-CA-C	6.54	120.59	111.56
1	D	202	ILE	N-CA-C	-5.96	104.54	110.62
1	A	110	LYS	N-CA-C	5.90	117.92	109.14
1	B	136	ASP	N-CA-C	-5.76	106.34	113.19
1	B	195	LYS	N-CA-C	-5.74	105.10	111.36
1	C	153	SER	N-CA-C	5.56	116.85	108.46
1	B	181	ASP	N-CA-C	5.49	122.50	110.80
1	B	202	ILE	N-CA-C	-5.49	105.17	110.72
1	A	203	SER	N-CA-C	5.24	117.07	111.36
1	A	166	SER	N-CA-C	-5.12	105.88	111.82
1	A	64	ASN	CB-CA-C	-5.07	104.73	111.63
1	A	154	ILE	CB-CA-C	-5.02	105.52	111.65
1	A	29	SER	CA-C-N	-5.02	114.61	122.65
1	A	29	SER	C-N-CA	-5.02	114.61	122.65

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	1599	0	1615	112	3
1	B	1630	0	1642	201	0
1	C	1631	0	1649	96	5
1	D	1607	0	1624	126	0
2	A	5	0	0	3	0
2	B	4	0	0	0	0
2	C	6	0	0	0	0
2	D	5	0	0	0	0
3	A	16	0	28	18	0
3	B	16	0	28	15	0
3	C	16	0	28	5	0
3	D	16	0	28	5	0
4	A	33	0	0	7	2
4	B	24	0	0	9	0
4	C	42	0	0	6	0
4	D	22	0	0	5	0
All	All	6672	0	6642	553	5

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 42.

All (553) 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:220:PHE:CD2	3:B:230:MPD:H51	1.16	1.69
1:A:7:LYS:HE2	1:A:7:LYS:N	1.12	1.45
1:B:220:PHE:HD2	3:B:230:MPD:C5	1.26	1.43
1:B:220:PHE:CD2	3:B:230:MPD:C5	1.99	1.34
1:B:96:HIS:HD2	1:B:119:ASP:OD1	1.10	1.33
1:D:182:MET:CE	1:D:182:MET:HA	1.56	1.33
1:A:7:LYS:N	1:A:7:LYS:CE	1.92	1.33
1:A:183:GLU:HG3	1:A:186:LYS:CB	1.60	1.31
1:B:13:ASP:CG	1:B:16:CYS:HB2	1.57	1.27
1:B:96:HIS:CD2	1:B:119:ASP:OD1	1.91	1.23
1:A:96:HIS:CE1	1:A:119:ASP:OD2	1.94	1.21

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:210:ALA:O	1:D:214:ILE:HD12	1.39	1.20
1:B:220:PHE:CE2	3:B:230:MPD:H51	1.77	1.20
1:B:216:PRO:HD2	1:B:219:SER:OG	1.43	1.18
1:A:183:GLU:CG	1:A:186:LYS:HB3	1.74	1.18
3:A:232:MPD:HM1	3:A:232:MPD:C5	1.70	1.17
1:B:181:ASP:HA	1:B:185:LEU:CD2	1.74	1.16
1:C:147:GLU:HG2	1:C:205:VAL:HG11	1.27	1.14
1:D:182:MET:CE	1:D:182:MET:CA	2.25	1.12
1:C:183:GLU:CB	1:C:184:PRO:HA	1.80	1.12
1:D:83:LYS:NZ	1:D:83:LYS:HB3	1.61	1.10
1:B:214:ILE:O	1:B:215:LEU:HD23	1.49	1.09
1:B:182:MET:O	1:B:186:LYS:HB2	1.51	1.09
1:B:181:ASP:HA	1:B:185:LEU:HD21	1.14	1.08
1:A:183:GLU:HG3	1:A:186:LYS:HB3	1.07	1.07
1:C:183:GLU:HB2	1:C:184:PRO:HA	1.13	1.06
1:D:134:ASN:ND2	1:D:138:GLY:O	1.91	1.04
1:D:182:MET:CA	1:D:182:MET:HE3	1.86	1.04
1:D:131:ASN:ND2	1:D:145:GLN:HE21	1.56	1.03
1:D:183:GLU:OE2	1:D:186:LYS:HB3	1.58	1.03
3:A:232:MPD:HM1	3:A:232:MPD:H52	1.40	1.03
1:C:183:GLU:HB2	1:C:184:PRO:CA	1.83	1.03
1:A:183:GLU:HA	1:A:185:LEU:N	1.72	1.03
1:D:159:ILE:HG21	3:D:232:MPD:H53	1.41	1.02
1:D:159:ILE:HG21	3:D:232:MPD:C5	1.89	1.01
3:C:233:MPD:O4	3:C:233:MPD:H12	1.57	1.01
1:B:181:ASP:CA	1:B:185:LEU:HD21	1.91	1.01
1:D:83:LYS:HB3	1:D:83:LYS:HZ3	1.20	1.00
1:B:183:GLU:HG3	1:B:184:PRO:HA	1.46	0.98
1:C:77:ASP:OD1	1:C:78:PHE:N	1.97	0.97
3:A:231:MPD:H52	3:A:231:MPD:O2	1.61	0.97
3:A:231:MPD:H53	3:A:231:MPD:H12	1.45	0.97
1:B:220:PHE:O	1:B:221:PHE:CG	2.17	0.97
1:A:84:ALA:O	1:A:85:LYS:HD2	1.64	0.96
1:B:220:PHE:CE2	4:B:250:HOH:O	2.19	0.96
1:D:210:ALA:O	1:D:214:ILE:CD1	2.13	0.96
1:B:64:ASN:OD1	4:B:254:HOH:O	1.83	0.95
1:D:147:GLU:CG	1:D:205:VAL:HG11	1.96	0.94
3:A:232:MPD:HM1	3:A:232:MPD:H53	1.47	0.94
1:A:75:ILE:HD13	1:A:91:THR:CG2	1.97	0.94
1:D:182:MET:HA	1:D:182:MET:HE3	0.94	0.93
1:D:182:MET:HE2	1:D:183:GLU:N	1.83	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:ILE:N	1:B:75:ILE:HD12	1.83	0.93
1:D:30:LYS:N	1:D:30:LYS:HD2	1.80	0.93
1:A:18:SER:O	1:A:22:GLU:HG3	1.69	0.93
1:A:183:GLU:HG3	1:A:186:LYS:HB2	1.50	0.92
1:B:116:TYR:HE2	3:B:231:MPD:CM	1.83	0.92
1:B:220:PHE:O	1:B:221:PHE:CD2	2.22	0.92
1:A:75:ILE:HD13	1:A:91:THR:HG22	1.51	0.91
1:B:158:LYS:HE2	1:B:158:LYS:HA	1.50	0.91
1:D:182:MET:HE2	1:D:182:MET:C	1.95	0.91
1:A:31:GLY:O	1:A:32:ILE:HG13	1.71	0.91
1:D:131:ASN:HD22	1:D:145:GLN:HE21	1.16	0.90
1:B:17:LEU:O	1:B:19:SER:N	2.05	0.90
1:B:10:LYS:O	1:B:12:GLY:N	2.05	0.89
1:B:18:SER:OG	1:B:78:PHE:HB2	1.71	0.89
1:D:182:MET:CA	1:D:182:MET:HE2	2.02	0.89
1:D:147:GLU:HG2	1:D:205:VAL:HG11	1.54	0.88
1:A:135:ASP:HB2	1:A:139:VAL:H	1.39	0.88
3:A:232:MPD:C5	3:A:232:MPD:CM	2.52	0.88
1:A:54:SER:OG	1:A:184:PRO:HB3	1.74	0.88
3:A:231:MPD:H12	3:A:231:MPD:C5	2.03	0.87
1:A:96:HIS:ND1	2:A:228:ZN:ZN	1.38	0.87
1:D:77:ASP:OD1	1:D:78:PHE:N	2.08	0.87
1:D:131:ASN:HB3	4:D:245:HOH:O	1.73	0.87
1:C:171:LYS:NZ	1:C:171:LYS:HB3	1.90	0.86
1:D:11:LEU:HD23	1:D:12:GLY:N	1.90	0.86
1:B:135:ASP:OD2	1:B:139:VAL:HB	1.76	0.85
1:C:171:LYS:HB3	1:C:171:LYS:HZ3	1.40	0.85
1:B:75:ILE:CG2	1:B:89:LEU:HD11	2.06	0.85
1:A:89:LEU:C	1:A:89:LEU:HD23	2.02	0.85
1:A:96:HIS:CE1	2:A:228:ZN:ZN	1.62	0.84
1:B:116:TYR:HE2	3:B:231:MPD:HM3	1.42	0.84
1:D:88:LEU:CD2	1:D:127:ARG:NH1	2.40	0.84
1:A:96:HIS:NE2	1:A:119:ASP:OD2	2.11	0.84
1:B:17:LEU:HD12	1:B:78:PHE:CZ	2.13	0.83
1:B:116:TYR:CE2	3:B:231:MPD:CM	2.61	0.83
1:A:183:GLU:HB3	1:A:187:THR:H	1.43	0.83
1:C:171:LYS:NZ	1:C:171:LYS:CB	2.38	0.82
1:B:75:ILE:HG23	1:B:89:LEU:HD11	1.62	0.82
1:B:38:TRP:HE3	1:B:203:SER:HB2	1.46	0.81
1:B:64:ASN:HA	4:B:254:HOH:O	1.80	0.81
1:B:38:TRP:CE3	1:B:203:SER:HB2	2.16	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:88:LEU:HD21	1:D:127:ARG:HH12	1.47	0.80
1:B:7:LYS:N	1:B:8:PRO:HD3	1.98	0.79
1:B:116:TYR:CE2	3:B:231:MPD:HM3	2.17	0.79
1:C:147:GLU:CG	1:C:205:VAL:HG11	2.09	0.79
1:C:214:ILE:HG22	1:C:215:LEU:HD13	1.64	0.79
1:B:13:ASP:OD1	1:B:16:CYS:HB2	1.81	0.78
1:B:218:SER:O	1:B:220:PHE:N	2.16	0.78
1:D:64:ASN:ND2	1:D:100:ASP:OD2	2.15	0.78
1:C:6:LEU:HD21	4:C:234:HOH:O	1.81	0.78
1:C:218:SER:HB2	4:C:273:HOH:O	1.83	0.77
1:A:17:LEU:HD23	1:A:78:PHE:CE1	2.19	0.77
1:B:80:MET:HE1	1:B:130:TYR:OH	1.83	0.77
1:C:11:LEU:HD23	1:C:11:LEU:C	2.09	0.77
1:C:88:LEU:CD1	1:C:127:ARG:HG2	2.15	0.77
1:C:183:GLU:CB	1:C:184:PRO:CA	2.53	0.77
1:D:8:PRO:HB2	1:D:222:GLU:CG	2.15	0.76
1:B:128:TYR:CD1	1:B:144:VAL:HG13	2.19	0.76
1:C:77:ASP:OD1	1:C:77:ASP:C	2.25	0.76
1:D:84:ALA:O	1:D:85:LYS:HB2	1.84	0.76
1:D:147:GLU:HG3	1:D:205:VAL:HG11	1.66	0.76
1:A:183:GLU:HA	1:A:184:PRO:C	2.11	0.76
1:B:214:ILE:C	1:B:215:LEU:HD23	2.10	0.76
1:D:92:LYS:HG2	1:D:123:ILE:HD13	1.66	0.76
1:A:75:ILE:CD1	1:A:91:THR:CG2	2.64	0.76
1:A:25:LEU:O	1:A:29:SER:HB3	1.87	0.75
1:B:218:SER:C	1:B:220:PHE:H	1.93	0.75
1:C:70:LEU:O	1:C:73:GLN:HG2	1.86	0.75
1:A:189:ARG:HB2	3:A:232:MPD:O4	1.86	0.75
1:A:31:GLY:O	1:A:32:ILE:CG1	2.35	0.75
1:D:75:ILE:N	1:D:75:ILE:HD12	2.02	0.75
1:B:183:GLU:CG	1:B:184:PRO:HA	2.17	0.74
1:A:89:LEU:HD22	1:A:126:VAL:HB	1.68	0.74
1:D:131:ASN:HD22	1:D:145:GLN:NE2	1.85	0.73
1:D:30:LYS:N	1:D:30:LYS:CD	2.49	0.73
1:D:88:LEU:HD22	1:D:127:ARG:NH1	2.03	0.73
1:B:22:GLU:OE1	1:B:75:ILE:N	2.21	0.73
1:A:96:HIS:HE1	4:A:254:HOH:O	1.71	0.73
1:A:135:ASP:OD2	1:A:139:VAL:HB	1.88	0.73
1:B:15:GLN:HA	1:B:15:GLN:NE2	2.04	0.73
1:D:88:LEU:HD22	1:D:127:ARG:HH11	1.54	0.73
1:A:135:ASP:HB3	1:A:137:ASN:H	1.53	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:162:SER:OG	1:A:164:ASP:OD1	2.05	0.72
1:B:220:PHE:CD2	3:B:230:MPD:H52	2.19	0.72
3:C:232:MPD:H12	3:C:232:MPD:O4	1.88	0.72
1:D:76:SER:O	1:D:77:ASP:CB	2.36	0.72
3:A:231:MPD:O2	3:A:231:MPD:C5	2.35	0.72
1:D:130:TYR:O	4:D:245:HOH:O	2.06	0.72
1:B:75:ILE:HG23	1:B:89:LEU:CD1	2.20	0.71
1:B:75:ILE:N	1:B:75:ILE:CD1	2.52	0.71
1:D:211:SER:HA	1:D:214:ILE:CD1	2.21	0.71
1:D:132:LEU:HB3	1:D:140:GLN:OE1	1.92	0.70
1:B:35:TYR:HB2	1:B:37:ILE:HD12	1.73	0.70
1:C:171:LYS:CB	1:C:171:LYS:HZ2	2.01	0.70
1:D:183:GLU:OE1	1:D:184:PRO:HA	1.91	0.70
1:B:77:ASP:OD1	1:B:78:PHE:N	2.24	0.70
1:C:218:SER:CB	4:C:273:HOH:O	2.40	0.70
1:B:220:PHE:HE2	4:B:250:HOH:O	1.60	0.70
1:D:45:VAL:HB	1:D:65:LEU:HB2	1.73	0.69
1:B:15:GLN:CD	1:B:15:GLN:O	2.35	0.69
1:B:75:ILE:CG2	1:B:89:LEU:CD1	2.70	0.69
1:B:158:LYS:HA	1:B:158:LYS:CE	2.23	0.69
1:D:182:MET:CE	1:D:182:MET:C	2.61	0.69
1:B:164:ASP:OD1	1:B:165:LEU:N	2.23	0.69
1:D:83:LYS:HB3	1:D:83:LYS:HZ2	1.58	0.69
1:A:7:LYS:N	1:A:7:LYS:CD	2.56	0.69
1:B:135:ASP:HB2	1:B:137:ASN:HD22	1.57	0.69
1:D:210:ALA:C	1:D:214:ILE:HD12	2.17	0.68
1:A:76:SER:O	1:A:77:ASP:HB2	1.92	0.68
1:D:115:LEU:HB2	1:D:162:SER:HB3	1.74	0.68
1:B:7:LYS:O	1:B:7:LYS:HG3	1.94	0.68
1:A:47:SER:HB2	4:A:259:HOH:O	1.94	0.67
1:A:89:LEU:C	1:A:89:LEU:CD2	2.67	0.67
1:B:11:LEU:O	1:B:11:LEU:HD23	1.94	0.67
1:B:7:LYS:O	1:B:7:LYS:CG	2.41	0.67
1:B:7:LYS:N	1:B:8:PRO:CD	2.57	0.67
1:B:154:ILE:HD12	1:B:157:PRO:HG2	1.76	0.67
1:C:88:LEU:HD12	1:C:127:ARG:HG2	1.76	0.67
1:D:17:LEU:HD21	1:D:221:PHE:CE1	2.30	0.67
1:D:61:ARG:HD2	1:D:104:GLU:OE1	1.95	0.67
1:A:89:LEU:HD23	1:A:89:LEU:O	1.95	0.66
1:B:38:TRP:HZ3	1:B:203:SER:HB3	1.60	0.66
1:B:38:TRP:CZ3	1:B:203:SER:HB3	2.30	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:89:LEU:C	1:C:89:LEU:HD12	2.20	0.66
1:B:52:ALA:HA	1:B:188:LEU:HD13	1.78	0.66
1:D:182:MET:HE2	1:D:183:GLU:H	1.61	0.66
1:D:131:ASN:ND2	1:D:145:GLN:NE2	2.39	0.65
1:A:84:ALA:O	1:A:85:LYS:CD	2.43	0.65
1:B:218:SER:C	1:B:220:PHE:N	2.51	0.65
1:B:220:PHE:O	1:B:221:PHE:CB	2.43	0.65
1:B:80:MET:CE	1:B:130:TYR:OH	2.44	0.65
1:A:10:LYS:HG3	1:A:10:LYS:O	1.94	0.65
1:B:75:ILE:HG21	1:B:89:LEU:HD11	1.79	0.65
1:B:22:GLU:CB	1:B:75:ILE:HD13	2.27	0.65
1:D:75:ILE:N	1:D:75:ILE:CD1	2.59	0.65
1:B:13:ASP:OD1	1:B:16:CYS:N	2.30	0.64
1:B:85:LYS:NZ	1:B:132:LEU:CD2	2.60	0.64
1:C:82:THR:HG22	1:C:130:TYR:OH	1.97	0.64
1:C:193:ILE:HD12	4:C:242:HOH:O	1.96	0.64
1:B:46:THR:HG22	1:B:47:SER:HB2	1.78	0.64
1:B:85:LYS:NZ	1:B:132:LEU:HD21	2.12	0.64
1:B:15:GLN:HA	1:B:15:GLN:HE21	1.62	0.64
1:D:92:LYS:CG	1:D:123:ILE:HD13	2.27	0.63
1:C:11:LEU:CB	1:C:223:ASN:OD1	2.46	0.63
1:D:88:LEU:CD1	1:D:127:ARG:NH1	2.62	0.63
1:D:83:LYS:NZ	1:D:83:LYS:CB	2.48	0.63
1:B:162:SER:OG	1:B:164:ASP:OD1	2.08	0.63
1:D:145:GLN:HB3	1:D:146:PRO:HD2	1.79	0.63
1:D:184:PRO:HD2	1:D:185:LEU:HD23	1.79	0.63
1:B:22:GLU:HB2	1:B:75:ILE:HD13	1.81	0.63
1:B:38:TRP:CE3	1:B:203:SER:CB	2.82	0.63
1:B:78:PHE:HE1	1:B:80:MET:HG3	1.63	0.63
1:D:42:PRO:HB3	1:D:68:THR:HG22	1.80	0.63
1:D:159:ILE:CG2	3:D:232:MPD:C5	2.74	0.63
1:C:88:LEU:HD13	1:C:127:ARG:HG2	1.79	0.62
1:A:109:SER:OG	4:A:235:HOH:O	2.15	0.62
3:A:232:MPD:H52	3:A:232:MPD:CM	2.22	0.62
1:B:77:ASP:HB3	1:B:90:LYS:HE3	1.81	0.62
1:B:220:PHE:C	1:B:221:PHE:CG	2.75	0.62
3:B:231:MPD:HO2	3:B:231:MPD:HO4	1.46	0.62
1:A:147:GLU:OE1	1:A:147:GLU:N	2.32	0.62
1:B:14:MET:HE3	1:B:79:GLN:HA	1.81	0.62
1:D:8:PRO:HB2	1:D:222:GLU:HG3	1.82	0.61
1:C:35:TYR:HB3	1:C:214:ILE:HD11	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:25:LEU:O	1:B:26:GLU:C	2.42	0.61
1:A:96:HIS:HD1	2:A:228:ZN:ZN	1.08	0.61
1:B:183:GLU:OE2	1:B:187:THR:HG23	2.01	0.61
1:B:17:LEU:O	1:B:18:SER:C	2.44	0.61
1:C:11:LEU:HB2	1:C:223:ASN:OD1	2.01	0.61
1:A:11:LEU:HD23	1:A:12:GLY:N	2.15	0.61
1:A:17:LEU:CD1	1:A:17:LEU:N	2.64	0.61
1:B:183:GLU:OE2	1:B:186:LYS:HB3	2.01	0.61
1:C:137:ASN:HD22	1:C:137:ASN:C	2.08	0.60
1:B:50:VAL:HG22	1:B:195:LYS:HG3	1.84	0.60
3:B:231:MPD:O4	3:B:231:MPD:O2	2.01	0.60
3:A:231:MPD:C5	3:A:231:MPD:C1	2.68	0.60
1:D:88:LEU:HD13	1:D:127:ARG:HG2	1.83	0.60
1:B:174:GLY:HA2	4:B:240:HOH:O	2.00	0.60
1:A:86:THR:HA	1:A:129:GLY:HA2	1.83	0.60
1:B:76:SER:O	1:B:77:ASP:CB	2.50	0.60
1:B:17:LEU:O	1:B:20:ALA:N	2.34	0.60
1:C:161:LEU:HD11	3:C:233:MPD:H53	1.83	0.60
1:B:47:SER:HB3	4:D:241:HOH:O	2.00	0.59
1:A:92:LYS:HA	1:A:122:VAL:O	2.02	0.59
1:C:147:GLU:HG2	1:C:205:VAL:CG1	2.18	0.59
1:A:183:GLU:CA	1:A:184:PRO:C	2.75	0.59
1:A:69:GLY:O	4:A:244:HOH:O	2.16	0.59
1:D:23:GLN:O	1:D:27:LYS:HG2	2.03	0.59
1:D:137:ASN:CG	1:D:137:ASN:O	2.46	0.59
1:C:171:LYS:HZ2	1:C:171:LYS:HB2	1.66	0.59
1:D:118:ALA:CB	1:D:159:ILE:HD13	2.33	0.59
1:C:84:ALA:O	1:C:85:LYS:HB2	2.02	0.58
1:A:183:GLU:HG2	1:A:186:LYS:HB3	1.76	0.58
1:D:76:SER:O	1:D:77:ASP:HB2	2.03	0.58
1:B:128:TYR:CE1	1:B:144:VAL:HG13	2.38	0.58
1:B:154:ILE:HD12	1:B:157:PRO:CG	2.34	0.58
1:B:154:ILE:CG1	4:B:243:HOH:O	2.51	0.58
1:C:10:LYS:CD	1:C:223:ASN:ND2	2.67	0.58
1:B:13:ASP:CB	1:B:16:CYS:HB2	2.34	0.58
1:B:15:GLN:NE2	1:B:15:GLN:CA	2.67	0.58
1:D:211:SER:HA	1:D:214:ILE:HD12	1.84	0.58
1:B:134:ASN:HA	1:B:139:VAL:O	2.05	0.57
1:B:135:ASP:HB2	1:B:137:ASN:ND2	2.20	0.57
1:A:214:ILE:CG2	1:A:214:ILE:O	2.53	0.57
1:D:42:PRO:HB3	1:D:68:THR:CG2	2.34	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:GLU:CD	1:A:183:GLU:N	2.63	0.56
1:B:216:PRO:HD2	1:B:219:SER:HG	1.66	0.56
1:B:216:PRO:CD	1:B:219:SER:OG	2.36	0.56
1:A:146:PRO:HG3	1:C:53:PRO:O	2.04	0.56
1:A:70:LEU:HB2	4:A:245:HOH:O	2.06	0.56
1:A:108:GLN:O	1:A:109:SER:HB2	2.05	0.56
1:D:128:TYR:HB3	1:D:147:GLU:HA	1.88	0.56
1:D:211:SER:HA	1:D:214:ILE:HD13	1.86	0.56
1:A:13:ASP:O	1:A:17:LEU:HD13	2.06	0.56
1:D:52:ALA:HA	1:D:188:LEU:HD13	1.87	0.56
1:B:85:LYS:HZ3	1:B:132:LEU:CD2	2.19	0.56
1:B:7:LYS:O	1:B:7:LYS:CD	2.54	0.56
1:D:210:ALA:C	1:D:214:ILE:CD1	2.76	0.56
1:A:22:GLU:HG2	1:A:75:ILE:HG12	1.88	0.55
1:B:38:TRP:CZ3	1:B:203:SER:CB	2.89	0.55
1:C:5:LEU:HD23	1:C:215:LEU:HD22	1.89	0.55
1:C:70:LEU:HD11	1:C:200:CYS:HB3	1.88	0.55
1:A:9:CYS:SG	1:A:17:LEU:HD12	2.47	0.55
1:C:10:LYS:HD3	1:C:223:ASN:ND2	2.21	0.55
1:C:128:TYR:CD1	1:C:144:VAL:HG13	2.41	0.55
1:C:133:LYS:HG2	1:C:143:GLU:HB2	1.87	0.55
1:C:18:SER:OG	1:C:78:PHE:HB2	2.07	0.55
1:C:61:ARG:HD2	1:C:104:GLU:OE1	2.07	0.55
1:A:132:LEU:N	1:A:132:LEU:HD12	2.21	0.55
1:A:163:SER:O	1:A:167:SER:HB3	2.07	0.55
1:C:13:ASP:OD2	1:C:16:CYS:HB2	2.07	0.55
1:C:28:THR:HG22	1:C:32:ILE:HD11	1.88	0.55
1:A:22:GLU:OE2	1:A:75:ILE:N	2.35	0.54
1:A:214:ILE:O	1:A:214:ILE:HG22	2.05	0.54
1:B:22:GLU:CA	1:B:75:ILE:HD13	2.38	0.54
1:D:12:GLY:HA2	1:D:14:MET:HE2	1.89	0.54
1:B:85:LYS:HZ1	1:B:132:LEU:HD21	1.72	0.54
1:D:192:ALA:O	1:D:196:ILE:HG13	2.07	0.54
1:B:22:GLU:HA	1:B:75:ILE:HD13	1.90	0.54
1:B:133:LYS:CB	1:B:133:LYS:NZ	2.70	0.54
1:A:89:LEU:CD2	1:A:89:LEU:O	2.55	0.54
1:B:63:LYS:O	1:B:64:ASN:HB2	2.08	0.54
1:C:45:VAL:HB	1:C:65:LEU:HB2	1.89	0.54
1:C:81:ASP:OD1	1:C:81:ASP:C	2.51	0.54
3:A:232:MPD:H53	3:A:232:MPD:CM	2.23	0.53
1:B:65:LEU:HD23	1:B:99:GLY:HA3	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:9:CYS:O	1:D:221:PHE:HA	2.08	0.53
1:B:28:THR:O	1:B:31:GLY:N	2.33	0.53
1:D:88:LEU:CD2	1:D:127:ARG:HH12	2.07	0.53
1:A:76:SER:O	1:A:77:ASP:CB	2.56	0.53
1:B:79:GLN:HB2	1:B:88:LEU:HB3	1.90	0.53
1:B:147:GLU:HG2	1:B:205:VAL:HG11	1.91	0.53
1:B:54:SER:O	1:B:55:ASP:HB2	2.09	0.52
1:C:119:ASP:O	1:C:157:PRO:HA	2.09	0.52
1:D:92:LYS:CG	1:D:123:ILE:CD1	2.87	0.52
1:C:129:GLY:HA3	1:C:145:GLN:NE2	2.23	0.52
1:C:133:LYS:HD2	1:C:143:GLU:OE2	2.08	0.52
1:D:73:GLN:HE22	1:D:201:TYR:HE1	1.56	0.52
1:A:130:TYR:CD1	1:A:130:TYR:C	2.87	0.52
1:A:185:LEU:O	1:A:186:LYS:C	2.52	0.52
1:B:135:ASP:C	1:B:137:ASN:H	2.17	0.52
1:D:9:CYS:SG	1:D:16:CYS:C	2.92	0.52
1:B:16:CYS:O	1:B:17:LEU:C	2.52	0.52
1:B:21:THR:O	1:B:22:GLU:C	2.52	0.52
1:D:147:GLU:HG2	1:D:205:VAL:CG1	2.32	0.52
1:B:22:GLU:HA	1:B:75:ILE:CD1	2.40	0.52
1:B:24:PHE:CZ	1:B:211:SER:HB3	2.45	0.52
1:B:172:ASP:C	1:B:174:GLY:H	2.17	0.52
1:A:11:LEU:HD23	1:A:11:LEU:C	2.35	0.52
1:A:131:ASN:C	1:A:132:LEU:HD12	2.34	0.52
1:B:220:PHE:CE2	3:B:230:MPD:C5	2.63	0.52
1:A:17:LEU:N	1:A:17:LEU:HD12	2.25	0.52
1:B:75:ILE:HD12	1:B:75:ILE:H	1.69	0.52
1:D:118:ALA:HB2	1:D:159:ILE:HD13	1.92	0.52
1:D:147:GLU:HG2	1:D:205:VAL:HG21	1.92	0.51
1:A:13:ASP:OD1	1:A:13:ASP:C	2.52	0.51
1:B:96:HIS:HD2	1:B:119:ASP:CG	2.03	0.51
1:B:137:ASN:ND2	1:B:137:ASN:C	2.67	0.51
1:B:17:LEU:C	1:B:19:SER:N	2.68	0.51
1:B:73:GLN:NE2	1:B:201:TYR:CE1	2.78	0.51
1:D:84:ALA:O	1:D:85:LYS:CB	2.55	0.51
1:B:76:SER:O	1:B:77:ASP:HB2	2.11	0.51
1:B:55:ASP:OD1	1:B:55:ASP:C	2.51	0.51
1:B:133:LYS:CB	1:B:133:LYS:HZ3	2.23	0.51
1:C:17:LEU:HD21	1:C:80:MET:CE	2.40	0.51
1:C:133:LYS:HD2	1:C:143:GLU:CD	2.36	0.51
1:B:183:GLU:HG3	1:B:184:PRO:CA	2.30	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:11:LEU:HD23	1:C:12:GLY:N	2.26	0.51
1:D:147:GLU:CD	1:D:147:GLU:H	2.18	0.51
1:B:47:SER:CB	4:D:241:HOH:O	2.57	0.51
1:C:137:ASN:ND2	1:C:139:VAL:H	2.09	0.51
1:A:54:SER:O	1:A:55:ASP:HB2	2.10	0.50
1:B:134:ASN:C	1:B:135:ASP:O	2.53	0.50
1:D:11:LEU:HD23	1:D:11:LEU:C	2.34	0.50
1:A:17:LEU:HD13	1:A:17:LEU:H	1.77	0.50
1:D:211:SER:CA	1:D:214:ILE:HD12	2.40	0.50
1:A:23:GLN:O	1:A:27:LYS:HG2	2.11	0.50
1:B:85:LYS:HZ1	1:B:132:LEU:CD2	2.24	0.50
1:D:97:ILE:O	1:D:117:THR:HA	2.11	0.50
1:D:73:GLN:NE2	1:D:201:TYR:OH	2.45	0.50
1:A:115:LEU:HB2	1:A:162:SER:HB3	1.93	0.50
1:B:116:TYR:HE2	3:B:231:MPD:HM1	1.72	0.50
1:A:28:THR:HB	1:A:37:ILE:HG21	1.93	0.50
1:D:28:THR:HB	1:D:37:ILE:HG21	1.93	0.50
1:C:17:LEU:CD2	1:C:80:MET:CE	2.90	0.50
1:A:142:PHE:CG	3:A:231:MPD:C1	2.95	0.50
1:B:26:GLU:HG2	1:B:27:LYS:HD3	1.92	0.49
1:B:113:THR:O	1:B:165:LEU:HG	2.11	0.49
1:B:158:LYS:HE2	1:B:158:LYS:CA	2.14	0.49
1:B:189:ARG:HG3	3:B:231:MPD:HO2	1.77	0.49
1:C:7:LYS:NZ	1:C:19:SER:HB3	2.27	0.49
1:A:31:GLY:C	1:A:32:ILE:HG13	2.35	0.49
1:D:76:SER:O	1:D:77:ASP:HB3	2.12	0.49
1:A:63:LYS:O	1:A:64:ASN:HB2	2.12	0.49
1:B:21:THR:O	1:B:23:GLN:N	2.45	0.49
1:B:23:GLN:HA	1:B:26:GLU:HB2	1.93	0.49
1:D:21:THR:O	1:D:22:GLU:C	2.55	0.49
1:D:88:LEU:HD13	1:D:127:ARG:NH1	2.27	0.49
1:B:140:GLN:O	1:B:217:ALA:N	2.31	0.49
1:B:21:THR:C	1:B:23:GLN:N	2.68	0.49
1:B:158:LYS:CE	1:B:158:LYS:CA	2.89	0.49
1:A:90:LYS:HA	1:A:124:GLY:O	2.12	0.49
1:A:7:LYS:N	1:A:7:LYS:NZ	2.58	0.48
1:B:125:ALA:HB2	1:B:152:GLU:OE2	2.13	0.48
1:B:133:LYS:HZ3	1:B:133:LYS:HB3	1.78	0.48
1:C:10:LYS:HD2	1:C:223:ASN:ND2	2.28	0.48
1:D:11:LEU:C	1:D:11:LEU:CD2	2.86	0.48
1:D:92:LYS:HG3	1:D:123:ILE:CD1	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:161:LEU:HD13	1:A:165:LEU:HD13	1.95	0.48
3:A:231:MPD:H52	3:A:231:MPD:C1	2.44	0.48
1:C:9:CYS:O	1:C:221:PHE:HA	2.13	0.48
1:B:135:ASP:C	1:B:137:ASN:N	2.71	0.48
1:B:189:ARG:HG3	3:B:231:MPD:O4	2.14	0.48
1:C:189:ARG:HA	3:C:233:MPD:H13	1.96	0.48
1:A:10:LYS:O	1:A:11:LEU:C	2.56	0.48
1:A:135:ASP:HB2	1:A:139:VAL:N	2.20	0.48
1:B:184:PRO:C	1:B:186:LYS:N	2.69	0.48
1:B:221:PHE:HD1	1:B:221:PHE:O	1.96	0.48
1:C:13:ASP:CG	1:C:16:CYS:HB2	2.39	0.48
1:D:94:ASP:OD1	1:D:121:ASN:ND2	2.46	0.48
1:D:145:GLN:HB3	1:D:146:PRO:CD	2.44	0.48
1:B:218:SER:O	1:B:221:PHE:N	2.39	0.48
1:A:17:LEU:CD1	1:A:17:LEU:H	2.25	0.48
1:A:29:SER:OG	1:A:71:LYS:HG2	2.14	0.48
1:B:22:GLU:OE1	1:B:74:GLN:HA	2.14	0.48
1:D:65:LEU:HA	1:D:65:LEU:HD23	1.57	0.48
1:B:75:ILE:CD1	1:B:75:ILE:H	2.23	0.47
1:C:129:GLY:HA3	1:C:145:GLN:HE22	1.79	0.47
3:C:233:MPD:O4	3:C:233:MPD:C1	2.30	0.47
1:B:17:LEU:HD12	1:B:78:PHE:CE1	2.49	0.47
1:C:15:GLN:HA	1:C:15:GLN:NE2	2.29	0.47
1:D:154:ILE:HD12	1:D:157:PRO:HG3	1.96	0.47
1:D:108:GLN:O	1:D:109:SER:HB2	2.13	0.47
1:B:158:LYS:C	1:B:159:ILE:HD12	2.39	0.47
1:C:7:LYS:HE3	1:C:23:GLN:OE1	2.15	0.47
1:C:76:SER:O	1:C:77:ASP:HB3	2.14	0.47
1:D:78:PHE:CD1	1:D:78:PHE:C	2.93	0.47
1:D:103:ILE:O	1:D:111:SER:HA	2.14	0.47
1:B:154:ILE:HG12	4:B:243:HOH:O	2.12	0.47
1:C:76:SER:O	1:C:77:ASP:CB	2.62	0.46
1:A:132:LEU:N	1:A:132:LEU:CD1	2.78	0.46
1:B:13:ASP:OD1	1:B:16:CYS:CB	2.60	0.46
1:B:11:LEU:O	1:B:11:LEU:CD2	2.63	0.46
1:A:75:ILE:CD1	1:A:91:THR:HG23	2.45	0.46
1:B:13:ASP:OD2	1:B:16:CYS:HB2	2.07	0.46
1:A:127:ARG:O	1:A:147:GLU:HB2	2.16	0.46
1:B:73:GLN:NE2	1:B:201:TYR:HE1	2.14	0.46
1:B:84:ALA:O	1:B:85:LYS:HB2	2.16	0.46
1:C:48:LEU:HD21	1:C:195:LYS:HB3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:LEU:HD23	1:A:140:GLN:OE1	2.15	0.46
1:B:132:LEU:O	1:B:133:LYS:HG2	2.16	0.46
1:C:90:LYS:NZ	1:C:90:LYS:HB3	2.30	0.46
1:C:11:LEU:C	1:C:11:LEU:CD2	2.82	0.45
1:D:9:CYS:O	1:D:222:GLU:N	2.49	0.45
1:B:70:LEU:O	1:B:73:GLN:HB3	2.15	0.45
1:C:27:LYS:HB3	1:C:27:LYS:HE2	1.67	0.45
1:B:88:LEU:HD12	1:B:89:LEU:H	1.81	0.45
1:C:60:ILE:HD12	1:C:60:ILE:N	2.31	0.45
1:D:8:PRO:HB2	1:D:222:GLU:HG2	1.92	0.45
1:D:41:ASP:HA	1:D:42:PRO:HA	1.70	0.45
1:D:66:ASN:C	1:D:67:ILE:HG13	2.41	0.45
1:D:158:LYS:HE2	1:D:158:LYS:HB2	1.64	0.45
1:A:142:PHE:CD1	3:A:231:MPD:H11	2.51	0.45
1:B:9:CYS:HB2	1:B:221:PHE:HB3	1.98	0.45
1:A:28:THR:HG22	1:A:32:ILE:CD1	2.47	0.45
1:A:81:ASP:OD2	1:A:84:ALA:HB3	2.17	0.45
1:C:17:LEU:CD2	1:C:80:MET:HE3	2.47	0.45
1:C:128:TYR:CE1	1:C:144:VAL:HG13	2.52	0.45
1:B:132:LEU:N	1:B:132:LEU:HD22	2.32	0.45
1:C:5:LEU:HD23	1:C:215:LEU:CD2	2.46	0.45
1:C:21:THR:O	1:C:25:LEU:HG	2.17	0.45
1:C:99:GLY:O	1:C:115:LEU:HD12	2.17	0.45
1:D:7:LYS:HA	1:D:8:PRO:HD3	1.76	0.45
1:D:131:ASN:HD21	1:D:145:GLN:HE21	1.57	0.45
1:A:36:ASP:O	1:A:206:HIS:HE1	2.00	0.45
1:B:135:ASP:OD2	1:B:139:VAL:CB	2.57	0.45
1:C:209:ARG:HD2	4:C:238:HOH:O	2.16	0.45
1:A:28:THR:HG22	1:A:32:ILE:HD11	1.99	0.44
1:B:123:ILE:HB	1:B:153:SER:HB3	1.98	0.44
1:A:27:LYS:HE3	1:A:27:LYS:HB3	1.56	0.44
1:A:41:ASP:HB3	1:A:71:LYS:HG3	2.00	0.44
1:B:18:SER:HB3	1:B:75:ILE:O	2.17	0.44
1:B:60:ILE:HD13	1:B:60:ILE:N	2.31	0.44
1:D:92:LYS:HG2	1:D:123:ILE:CD1	2.42	0.44
1:B:7:LYS:O	1:B:7:LYS:HD2	2.17	0.44
1:C:88:LEU:CD1	1:C:127:ARG:CG	2.89	0.44
1:D:169:LEU:O	1:D:170:GLU:C	2.56	0.44
1:A:74:GLN:O	1:A:91:THR:HG22	2.17	0.44
1:C:66:ASN:C	1:C:67:ILE:HG13	2.41	0.44
1:A:104:GLU:HG3	1:A:111:SER:OG	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:36:ASP:O	1:B:206:HIS:HE1	2.01	0.44
1:B:74:GLN:C	1:B:75:ILE:HD12	2.41	0.44
1:B:166:SER:HB3	4:B:242:HOH:O	2.17	0.44
1:B:218:SER:O	1:B:219:SER:C	2.61	0.44
1:C:11:LEU:HB3	1:C:223:ASN:OD1	2.18	0.44
1:C:88:LEU:HD13	1:C:127:ARG:CG	2.46	0.44
1:C:12:GLY:HA2	1:C:14:MET:HE2	2.00	0.44
1:C:193:ILE:CD1	4:C:242:HOH:O	2.59	0.44
1:C:108:GLN:O	1:C:109:SER:HB2	2.17	0.44
1:A:183:GLU:N	1:A:183:GLU:OE1	2.51	0.44
1:B:75:ILE:O	1:B:75:ILE:HG22	2.17	0.44
1:B:83:LYS:HB3	1:B:83:LYS:HE2	1.55	0.44
1:B:181:ASP:CB	1:B:185:LEU:HD21	2.46	0.44
1:A:183:GLU:CA	1:A:185:LEU:N	2.62	0.43
1:D:11:LEU:HD23	1:D:12:GLY:H	1.78	0.43
1:A:13:ASP:OD2	1:A:16:CYS:HB2	2.18	0.43
1:C:28:THR:HG22	1:C:32:ILE:CD1	2.47	0.43
1:D:77:ASP:CG	1:D:79:GLN:HE21	2.26	0.43
1:A:209:ARG:HG3	4:A:249:HOH:O	2.19	0.43
3:A:231:MPD:H53	3:A:231:MPD:C1	2.30	0.43
1:B:78:PHE:CE1	1:B:80:MET:HG3	2.49	0.43
1:B:85:LYS:HA	1:B:85:LYS:HD3	1.47	0.43
1:B:142:PHE:N	4:B:239:HOH:O	2.02	0.43
1:D:121:ASN:O	1:D:154:ILE:O	2.36	0.43
1:A:159:ILE:HG21	3:A:232:MPD:H53	1.99	0.43
1:C:156:GLU:HA	1:C:157:PRO:HD2	1.78	0.43
1:A:142:PHE:CD1	3:A:231:MPD:C1	3.02	0.43
1:C:60:ILE:N	1:C:60:ILE:CD1	2.81	0.43
1:A:67:ILE:HG12	1:A:97:ILE:HG12	2.01	0.43
1:C:15:GLN:CD	1:C:15:GLN:O	2.62	0.43
1:D:70:LEU:HD11	1:D:200:CYS:HB3	1.99	0.43
1:B:78:PHE:HE1	1:B:80:MET:CG	2.29	0.43
1:B:10:LYS:C	1:B:12:GLY:N	2.75	0.42
1:B:70:LEU:HD11	1:B:200:CYS:HB3	2.01	0.42
1:B:91:THR:HB	1:B:201:TYR:OH	2.19	0.42
1:B:14:MET:C	1:B:16:CYS:H	2.26	0.42
1:D:88:LEU:HD21	1:D:127:ARG:NH1	2.09	0.42
1:D:141:HIS:ND1	1:D:215:LEU:O	2.51	0.42
1:A:21:THR:O	1:A:22:GLU:C	2.60	0.42
1:B:14:MET:C	1:B:16:CYS:N	2.77	0.42
1:A:78:PHE:CD1	1:A:78:PHE:C	2.97	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:ILE:O	1:B:77:ASP:N	2.49	0.42
1:B:132:LEU:HG	1:B:140:GLN:NE2	2.35	0.42
1:D:186:LYS:O	1:D:186:LYS:HG3	2.20	0.42
1:C:137:ASN:C	1:C:137:ASN:ND2	2.74	0.42
1:D:28:THR:O	1:D:31:GLY:N	2.44	0.42
1:D:104:GLU:HG2	1:D:106:THR:HG23	2.02	0.42
1:A:145:GLN:HE21	1:A:145:GLN:HB2	1.63	0.42
1:B:164:ASP:CG	1:B:165:LEU:N	2.77	0.42
1:C:123:ILE:HB	1:C:153:SER:HB3	2.02	0.42
1:D:38:TRP:NE1	1:D:206:HIS:CE1	2.88	0.42
1:A:48:LEU:HD12	1:A:48:LEU:HA	1.87	0.41
1:A:96:HIS:CE1	4:A:254:HOH:O	2.57	0.41
1:B:21:THR:C	1:B:23:GLN:H	2.28	0.41
1:B:77:ASP:OD1	1:B:77:ASP:C	2.61	0.41
1:B:209:ARG:HG2	1:B:213:LYS:HE3	2.02	0.41
1:C:137:ASN:HD22	1:C:139:VAL:H	1.68	0.41
1:C:145:GLN:HE21	1:C:145:GLN:HB2	1.47	0.41
1:C:183:GLU:HB3	1:C:184:PRO:HA	1.89	0.41
1:D:188:LEU:HG	3:D:232:MPD:H13	2.01	0.41
1:A:49:ASP:O	1:A:195:LYS:NZ	2.53	0.41
1:B:29:SER:OG	1:B:71:LYS:HB2	2.19	0.41
1:B:82:THR:O	1:B:85:LYS:HE3	2.20	0.41
1:B:12:GLY:O	1:B:13:ASP:C	2.63	0.41
1:B:35:TYR:CB	1:B:37:ILE:HD12	2.47	0.41
1:B:171:LYS:HE2	1:B:171:LYS:HB2	1.72	0.41
1:C:103:ILE:O	1:C:111:SER:HA	2.20	0.41
1:D:205:VAL:O	1:D:206:HIS:C	2.64	0.41
1:B:92:LYS:HA	1:B:122:VAL:O	2.21	0.41
1:B:186:LYS:HE2	1:B:190:GLN:HE22	1.86	0.41
1:C:10:LYS:HD2	1:C:223:ASN:HD21	1.85	0.41
1:C:95:LEU:HA	1:C:95:LEU:HD23	1.73	0.41
1:D:128:TYR:CD1	1:D:144:VAL:HG13	2.55	0.41
1:D:11:LEU:CD2	1:D:12:GLY:N	2.74	0.41
3:D:232:MPD:O4	3:D:232:MPD:H12	2.20	0.41
1:A:208:ILE:HD13	1:A:208:ILE:HG21	1.81	0.40
1:B:16:CYS:O	1:B:17:LEU:O	2.38	0.40
1:C:128:TYR:HD1	1:C:129:GLY:O	2.04	0.40
1:D:41:ASP:CG	1:D:69:GLY:H	2.29	0.40
1:D:145:GLN:HB2	4:D:238:HOH:O	2.21	0.40
1:B:88:LEU:HD13	1:B:127:ARG:HG2	2.02	0.40
1:B:147:GLU:H	1:B:147:GLU:HG3	1.14	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:88:LEU:HD13	1:D:127:ARG:HH11	1.86	0.40
1:D:186:LYS:HB2	1:D:186:LYS:HE3	1.78	0.40
1:A:92:LYS:HG2	1:A:123:ILE:HG12	2.03	0.40
1:D:8:PRO:HG2	1:D:222:GLU:HG2	2.03	0.40
1:A:118:ALA:HA	1:A:158:LYS:O	2.22	0.40
1:B:172:ASP:O	1:B:174:GLY:N	2.52	0.40
1:C:137:ASN:HD22	1:C:138:GLY:N	2.20	0.40
1:B:97:ILE:HB	1:B:118:ALA:HB3	2.02	0.40
1:C:41:ASP:HA	1:C:42:PRO:HA	1.90	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:26:GLU:CG	4:A:254:HOH:O[1_455]	1.32	0.88
1:A:119:ASP:OD2	1:C:26:GLU:OE2[1_655]	1.59	0.61
1:A:96:HIS:CE1	1:C:26:GLU:OE2[1_655]	1.97	0.23
1:C:26:GLU:CD	4:A:254:HOH:O[1_455]	2.00	0.20
1:A:96:HIS:NE2	1:C:26:GLU:OE2[1_655]	2.05	0.15

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	204/227 (90%)	190 (93%)	8 (4%)	6 (3%)	3	6
1	B	208/227 (92%)	178 (86%)	20 (10%)	10 (5%)	2	2
1	C	208/227 (92%)	195 (94%)	10 (5%)	3 (1%)	9	19
1	D	205/227 (90%)	189 (92%)	12 (6%)	4 (2%)	6	12
All	All	825/908 (91%)	752 (91%)	50 (6%)	23 (3%)	4	6

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	55	ASP
1	A	77	ASP
1	A	135	ASP
1	B	11	LEU
1	B	17	LEU
1	B	18	SER
1	B	77	ASP
1	B	221	PHE
1	C	77	ASP
1	D	77	ASP
1	A	185	LEU
1	B	29	SER
1	B	173	SER
1	B	219	SER
1	D	11	LEU
1	B	12	GLY
1	B	55	ASP
1	C	219	SER
1	A	184	PRO
1	D	8	PRO
1	D	78	PHE
1	A	13	ASP
1	C	218	SER

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	182/198 (92%)	158 (87%)	24 (13%)	4	8
1	B	186/198 (94%)	156 (84%)	30 (16%)	2	4
1	C	186/198 (94%)	163 (88%)	23 (12%)	4	9
1	D	183/198 (92%)	158 (86%)	25 (14%)	3	7
All	All	737/792 (93%)	635 (86%)	102 (14%)	3	7

All (102) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LYS
1	A	17	LEU
1	A	19	SER
1	A	30	LYS
1	A	34	GLN
1	A	61	ARG
1	A	71	LYS
1	A	85	LYS
1	A	89	LEU
1	A	91	THR
1	A	106	THR
1	A	115	LEU
1	A	120	THR
1	A	144	VAL
1	A	156	GLU
1	A	159	ILE
1	A	160	THR
1	A	165	LEU
1	A	166	SER
1	A	183	GLU
1	A	189	ARG
1	A	203	SER
1	A	214	ILE
1	A	218	SER
1	B	10	LYS
1	B	15	GLN
1	B	17	LEU
1	B	47	SER
1	B	48	LEU
1	B	60	ILE
1	B	61	ARG
1	B	66	ASN
1	B	75	ILE
1	B	83	LYS
1	B	85	LYS
1	B	92	LYS
1	B	115	LEU
1	B	117	THR
1	B	121	ASN
1	B	133	LYS
1	B	137	ASN
1	B	139	VAL
1	B	140	GLN

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Mol	Chain	Res	Type
1	B	147	GLU
1	B	148	THR
1	B	158	LYS
1	B	165	LEU
1	B	171	LYS
1	B	185	LEU
1	B	189	ARG
1	B	190	GLN
1	B	195	LYS
1	B	221	PHE
1	B	222	GLU
1	C	6	LEU
1	C	17	LEU
1	C	34	GLN
1	C	60	ILE
1	C	61	ARG
1	C	77	ASP
1	C	90	LYS
1	C	121	ASN
1	C	137	ASN
1	C	147	GLU
1	C	158	LYS
1	C	165	LEU
1	C	166	SER
1	C	171	LYS
1	C	183	GLU
1	C	185	LEU
1	C	190	GLN
1	C	193	ILE
1	C	195	LYS
1	C	203	SER
1	C	213	LYS
1	C	215	LEU
1	C	219	SER
1	D	10	LYS
1	D	23	GLN
1	D	30	LYS
1	D	74	GLN
1	D	75	ILE
1	D	83	LYS
1	D	89	LEU
1	D	92	LYS

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Mol	Chain	Res	Type
1	D	94	ASP
1	D	110	LYS
1	D	119	ASP
1	D	131	ASN
1	D	136	ASP
1	D	147	GLU
1	D	158	LYS
1	D	165	LEU
1	D	166	SER
1	D	167	SER
1	D	182	MET
1	D	185	LEU
1	D	187	THR
1	D	195	LYS
1	D	203	SER
1	D	219	SER
1	D	222	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	GLN
1	A	79	GLN
1	A	96	HIS
1	A	145	GLN
1	B	15	GLN
1	B	66	ASN
1	B	72	ASN
1	B	74	GLN
1	B	96	HIS
1	B	108	GLN
1	B	137	ASN
1	B	140	GLN
1	B	190	GLN
1	C	15	GLN
1	C	34	GLN
1	C	66	ASN
1	C	73	GLN
1	C	79	GLN
1	C	131	ASN
1	C	137	ASN
1	C	145	GLN

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Mol	Chain	Res	Type
1	C	190	GLN
1	D	23	GLN
1	D	34	GLN
1	D	66	ASN
1	D	73	GLN
1	D	79	GLN
1	D	108	GLN
1	D	121	ASN
1	D	131	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 28 ligands modelled in this entry, 20 are monoatomic - leaving 8 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	MPD	D	231	-	7,7,7	0.37	0	9,10,10	0.61	0
3	MPD	C	232	-	7,7,7	0.43	0	9,10,10	0.66	0
3	MPD	A	232	-	7,7,7	0.44	0	9,10,10	0.64	0
3	MPD	A	231	-	7,7,7	0.30	0	9,10,10	0.23	0
3	MPD	D	232	-	7,7,7	0.38	0	9,10,10	0.46	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	MPD	B	230	-	7,7,7	0.34	0	9,10,10	0.31	0
3	MPD	C	233	-	7,7,7	0.45	0	9,10,10	0.80	0
3	MPD	B	231	-	7,7,7	0.35	0	9,10,10	0.35	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MPD	D	231	-	-	0/5/5/5	-
3	MPD	C	232	-	-	3/5/5/5	-
3	MPD	A	232	-	-	1/5/5/5	-
3	MPD	A	231	-	-	3/5/5/5	-
3	MPD	D	232	-	-	3/5/5/5	-
3	MPD	B	230	-	-	3/5/5/5	-
3	MPD	C	233	-	-	2/5/5/5	-
3	MPD	B	231	-	-	3/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	231	MPD	C2-C3-C4-O4
3	A	231	MPD	C2-C3-C4-C5
3	A	232	MPD	C2-C3-C4-C5
3	B	230	MPD	C1-C2-C3-C4
3	B	230	MPD	C2-C3-C4-C5
3	C	232	MPD	C1-C2-C3-C4
3	D	232	MPD	C2-C3-C4-C5
3	C	232	MPD	O2-C2-C3-C4
3	B	231	MPD	C1-C2-C3-C4
3	C	232	MPD	CM-C2-C3-C4
3	C	233	MPD	C1-C2-C3-C4
3	D	232	MPD	O2-C2-C3-C4
3	B	231	MPD	C2-C3-C4-O4
3	D	232	MPD	C2-C3-C4-O4

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Mol	Chain	Res	Type	Atoms
3	A	231	MPD	C1-C2-C3-C4
3	B	230	MPD	CM-C2-C3-C4
3	B	231	MPD	CM-C2-C3-C4
3	C	233	MPD	CM-C2-C3-C4

There are no ring outliers.

7 monomers are involved in 43 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	232	MPD	1	0
3	A	232	MPD	8	0
3	A	231	MPD	10	0
3	D	232	MPD	5	0
3	B	230	MPD	6	0
3	C	233	MPD	4	0
3	B	231	MPD	9	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, 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	208/227 (91%)	0.19	3 (1%) 73 69	28, 44, 76, 89	0
1	B	212/227 (93%)	0.70	17 (8%) 18 14	27, 59, 98, 110	0
1	C	212/227 (93%)	0.12	5 (2%) 59 54	26, 42, 69, 77	0
1	D	209/227 (92%)	0.49	14 (6%) 24 19	29, 50, 83, 96	0
All	All	841/908 (92%)	0.37	39 (4%) 37 32	26, 48, 82, 110	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	11	LEU	5.4
1	B	221	PHE	5.0
1	B	180	PRO	4.9
1	B	222	GLU	4.3
1	C	5	LEU	4.2
1	D	7	LYS	4.1
1	A	221	PHE	4.1
1	B	8	PRO	4.0
1	D	9	CYS	4.0
1	B	7	LYS	3.9
1	D	182	MET	3.9
1	A	11	LEU	3.7
1	B	10	LYS	3.3
1	B	182	MET	3.1
1	A	222	GLU	3.1
1	C	76	SER	3.1
1	D	184	PRO	3.0
1	D	185	LEU	2.9
1	B	9	CYS	2.8
1	D	174	GLY	2.8
1	D	222	GLU	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	184	PRO	2.7
1	D	23	GLN	2.6
1	B	18	SER	2.6
1	C	184	PRO	2.6
1	D	12	GLY	2.5
1	D	11	LEU	2.5
1	C	183	GLU	2.4
1	B	27	LYS	2.4
1	B	136	ASP	2.4
1	B	19	SER	2.4
1	D	76	SER	2.4
1	C	7	LYS	2.4
1	D	75	ILE	2.3
1	B	14	MET	2.3
1	D	16	CYS	2.1
1	B	17	LEU	2.0
1	B	175	ASN	2.0
1	D	8	PRO	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 oligosaccharides 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. 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	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ZN	C	231	1/1	-0.50	0.71	89,89,89,89	1
2	ZN	A	230	1/1	0.62	0.44	77,77,77,77	1
2	ZN	D	228	1/1	0.71	0.19	67,67,67,67	1
2	ZN	B	229	1/1	0.73	0.37	84,84,84,84	1
2	ZN	D	230	1/1	0.73	0.19	70,70,70,70	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MPD	B	230	8/8	0.85	0.18	73,73,74,74	0
2	ZN	A	228	1/1	0.88	0.10	75,75,75,75	1
2	ZN	C	228	1/1	0.89	0.10	47,47,47,47	1
2	ZN	B	228	1/1	0.91	0.09	49,49,49,49	1
3	MPD	B	231	8/8	0.91	0.18	67,69,71,72	0
3	MPD	A	231	8/8	0.92	0.19	73,74,75,77	0
2	ZN	D	227	1/1	0.93	0.14	82,82,82,82	0
3	MPD	D	231	8/8	0.93	0.16	61,62,67,68	0
3	MPD	D	232	8/8	0.93	0.14	52,54,58,61	0
2	ZN	C	229	1/1	0.94	0.37	80,80,80,80	1
2	ZN	B	227	1/1	0.94	0.09	81,81,81,81	0
3	MPD	C	232	8/8	0.94	0.12	47,49,51,53	0
2	ZN	A	229	1/1	0.94	0.15	43,43,43,43	1
3	MPD	A	232	8/8	0.94	0.13	42,45,52,52	0
2	ZN	D	229	1/1	0.96	0.21	52,52,52,52	1
3	MPD	C	233	8/8	0.96	0.11	43,46,53,53	0
2	ZN	C	230	1/1	0.97	0.08	38,38,38,38	1
2	ZN	B	226	1/1	0.98	0.04	47,47,47,47	0
2	ZN	C	226	1/1	0.98	0.05	34,34,34,34	0
2	ZN	D	226	1/1	0.98	0.07	38,38,38,38	0
2	ZN	C	227	1/1	0.99	0.04	41,41,41,41	0
2	ZN	A	227	1/1	0.99	0.02	37,37,37,37	0
2	ZN	A	226	1/1	1.00	0.03	35,35,35,35	0

6.5 Other polymers

There are no such residues in this entry.