



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:29 PM UTC

PDB ID : 6W8S / pdb_00006w8s
Title : Crystal structure of metacaspase 4 from Arabidopsis
Authors : Zhu, P.; Yu, X.H.; Wang, C.; Zhang, Q.; Liu, W.; McSweeney, S.; Shanklin, J.; Lam, E.; Liu, Q.
Deposited on : 2020-03-21
Resolution : 3.48 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

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

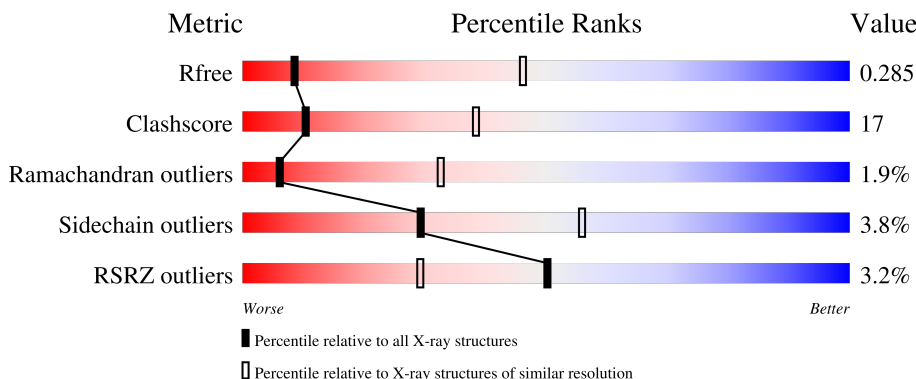
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.48 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}	180053	1083 (3.52-3.44)
Clashscore	190562	1139 (3.52-3.44)
Ramachandran outliers	187476	1111 (3.52-3.44)
Sidechain outliers	187428	1112 (3.52-3.44)
RSRZ outliers	180081	1082 (3.52-3.44)

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	426	4% 54% 29% • 15%
1	B	426	2% 54% 28% • 16%
1	C	426	3% 53% 29% • 15%
1	D	426	2% 52% 30% • 16%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10908 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 Metacaspase-4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	362	Total	C	N	O	S	0	0	0
			2728	1684	465	561	18			
1	B	357	Total	C	N	O	S	0	0	0
			2698	1666	461	553	18			
1	C	361	Total	C	N	O	S	0	0	0
			2732	1688	465	561	18			
1	D	358	Total	C	N	O	S	0	0	0
			2710	1674	462	556	18			

There are 32 discrepancies between the modelled and reference sequences:

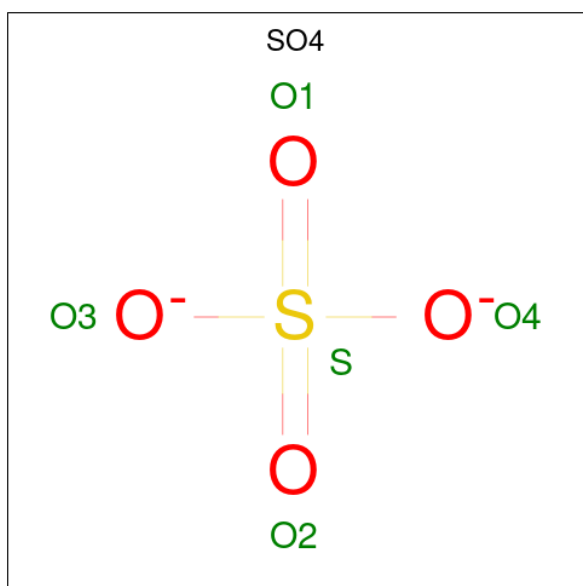
Chain	Residue	Modelled	Actual	Comment	Reference
A	419	VAL	-	expression tag	UNP O64517
A	420	GLU	-	expression tag	UNP O64517
A	421	HIS	-	expression tag	UNP O64517
A	422	HIS	-	expression tag	UNP O64517
A	423	HIS	-	expression tag	UNP O64517
A	424	HIS	-	expression tag	UNP O64517
A	425	HIS	-	expression tag	UNP O64517
A	426	HIS	-	expression tag	UNP O64517
B	419	VAL	-	expression tag	UNP O64517
B	420	GLU	-	expression tag	UNP O64517
B	421	HIS	-	expression tag	UNP O64517
B	422	HIS	-	expression tag	UNP O64517
B	423	HIS	-	expression tag	UNP O64517
B	424	HIS	-	expression tag	UNP O64517
B	425	HIS	-	expression tag	UNP O64517
B	426	HIS	-	expression tag	UNP O64517
C	419	VAL	-	expression tag	UNP O64517
C	420	GLU	-	expression tag	UNP O64517
C	421	HIS	-	expression tag	UNP O64517
C	422	HIS	-	expression tag	UNP O64517
C	423	HIS	-	expression tag	UNP O64517

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Chain	Residue	Modelled	Actual	Comment	Reference
C	424	HIS	-	expression tag	UNP O64517
C	425	HIS	-	expression tag	UNP O64517
C	426	HIS	-	expression tag	UNP O64517
D	419	VAL	-	expression tag	UNP O64517
D	420	GLU	-	expression tag	UNP O64517
D	421	HIS	-	expression tag	UNP O64517
D	422	HIS	-	expression tag	UNP O64517
D	423	HIS	-	expression tag	UNP O64517
D	424	HIS	-	expression tag	UNP O64517
D	425	HIS	-	expression tag	UNP O64517
D	426	HIS	-	expression tag	UNP O64517

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

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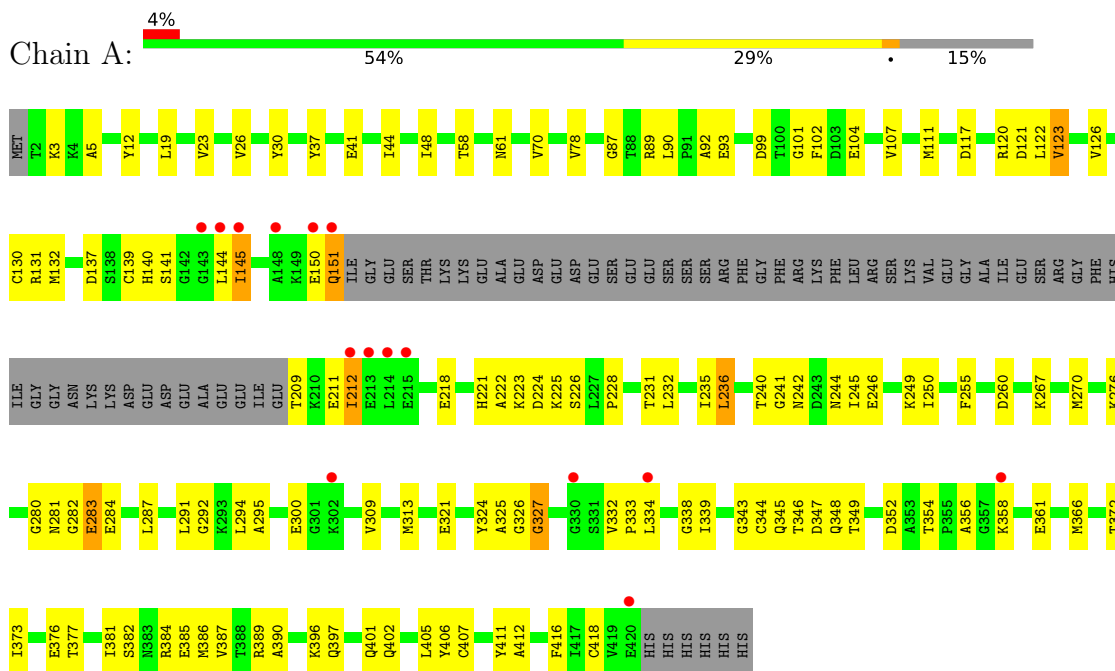
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

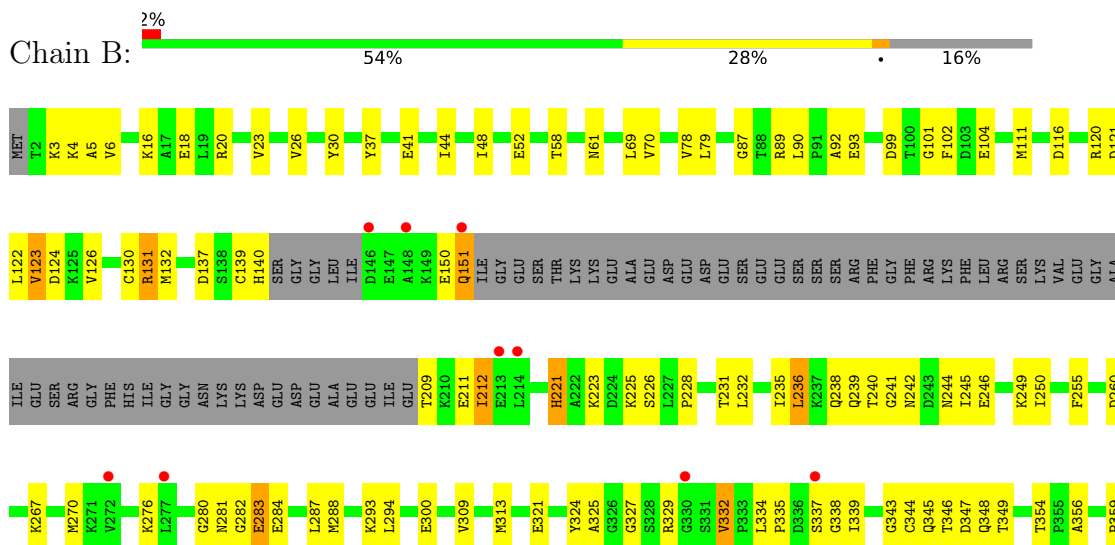
3 Residue-property plots

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: Metacaspase-4



• Molecule 1: Metacaspase-4



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	125.91Å 284.44Å 57.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.80 – 3.48 38.80 – 3.48	Depositor EDS
% Data completeness (in resolution range)	88.9 (38.80-3.48) 88.7 (38.80-3.48)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.43 (at 3.48Å)	Xtriage
Refinement program	PHENIX 1.10.1 _2155	Depositor
R, R_{free}	0.250 , 0.284 0.254 , 0.285	Depositor DCC
R_{free} test set	1242 reflections (4.47%)	wwPDB-VP
Wilson B-factor (Å ²)	89.7	Xtriage
Anisotropy	0.058	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 80.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	10908	wwPDB-VP
Average B, all atoms (Å ²)	93.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.71% of the height of the origin peak. No significant pseudotranslation is detected.*

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: 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	0.34	0/2766	0.85	2/3732 (0.1%)
1	B	0.29	0/2735	0.86	8/3688 (0.2%)
1	C	0.28	0/2769	0.86	4/3734 (0.1%)
1	D	0.36	0/2747	0.90	6/3704 (0.2%)
All	All	0.32	0/11017	0.87	20/14858 (0.1%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	217	GLY	N-CA-C	9.09	128.10	115.00
1	B	329	ARG	CA-C-N	7.14	133.28	120.79
1	B	329	ARG	C-N-CA	7.14	133.28	120.79
1	D	214	LEU	CA-C-N	-6.82	110.28	122.15
1	D	214	LEU	C-N-CA	-6.82	110.28	122.15
1	C	220	ILE	N-CA-C	6.74	117.54	108.11
1	D	332	VAL	CA-C-N	6.62	126.66	119.90
1	D	332	VAL	C-N-CA	6.62	126.66	119.90
1	C	332	VAL	N-CA-C	6.42	114.96	107.77
1	D	220	ILE	N-CA-C	6.32	116.95	108.11
1	A	326	GLY	N-CA-C	-6.10	107.35	115.40
1	B	338	GLY	N-CA-C	-5.81	99.42	113.18
1	C	338	GLY	N-CA-C	-5.52	100.09	113.18
1	B	329	ARG	N-CA-C	-5.43	107.30	114.31
1	A	338	GLY	N-CA-C	-5.43	100.31	113.18
1	B	337	SER	CA-C-N	5.41	132.02	121.41
1	B	337	SER	C-N-CA	5.41	132.02	121.41
1	B	332	VAL	CA-C-N	5.20	124.96	119.76
1	B	332	VAL	C-N-CA	5.20	124.96	119.76
1	D	380	GLU	N-CA-C	-5.10	101.39	109.76

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	2728	0	2672	91	0
1	B	2698	0	2645	84	0
1	C	2732	0	2679	99	0
1	D	2710	0	2660	97	0
2	A	15	0	0	0	0
2	B	5	0	0	0	0
2	C	15	0	0	1	0
2	D	5	0	0	1	0
All	All	10908	0	10656	367	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 17.

All (367) 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:20:ARG:HD3	1:B:359:PRO:HB3	1.39	1.04
1:C:20:ARG:HD3	1:C:359:PRO:HB3	1.44	0.98
1:D:66:LEU:HD11	1:D:122:LEU:HD13	1.51	0.91
1:A:37:TYR:HD1	1:A:418:CYS:HG	1.26	0.83
1:C:145:ILE:HG23	1:C:334:LEU:HD12	1.61	0.81
1:D:145:ILE:HG23	1:D:334:LEU:HD12	1.65	0.79
1:D:37:TYR:HD1	1:D:418:CYS:HG	1.30	0.77
1:A:348:GLN:NE2	1:A:402:GLN:H	1.83	0.76
1:A:145:ILE:HG21	1:A:406:TYR:HB3	1.66	0.76
1:C:348:GLN:NE2	1:C:402:GLN:H	1.83	0.76
1:D:348:GLN:NE2	1:D:402:GLN:H	1.85	0.75
1:C:145:ILE:HG21	1:C:406:TYR:HB3	1.69	0.75
1:B:348:GLN:NE2	1:B:402:GLN:H	1.84	0.74
1:D:145:ILE:HG21	1:D:406:TYR:HB3	1.68	0.74
1:D:49:ASP:OD1	1:D:50:THR:N	2.22	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:345:GLN:H	1:B:348:GLN:NE2	1.88	0.71
1:D:13:PRO:HD3	1:D:49:ASP:OD2	1.90	0.71
1:A:145:ILE:HG23	1:A:334:LEU:HD12	1.73	0.70
1:D:345:GLN:H	1:D:348:GLN:NE2	1.90	0.70
1:D:3:LYS:HG2	1:D:78:VAL:HB	1.74	0.70
1:C:345:GLN:H	1:C:348:GLN:NE2	1.90	0.69
1:C:70:VAL:HG13	1:C:126:VAL:HG12	1.74	0.69
1:B:3:LYS:HG2	1:B:78:VAL:HB	1.74	0.69
1:A:345:GLN:H	1:A:348:GLN:NE2	1.90	0.68
1:A:121:ASP:OD1	1:A:332:VAL:HG21	1.92	0.68
1:A:339:ILE:HG13	1:A:407:CYS:HB3	1.76	0.68
1:B:339:ILE:HG13	1:B:407:CYS:HB3	1.76	0.67
1:A:120:ARG:HH11	1:A:144:LEU:HG	1.59	0.67
1:A:70:VAL:HG13	1:A:126:VAL:HG12	1.76	0.67
1:C:3:LYS:HG2	1:C:78:VAL:HB	1.76	0.67
1:B:70:VAL:HG13	1:B:126:VAL:HG12	1.78	0.66
1:A:120:ARG:NE	1:A:332:VAL:O	2.25	0.65
1:D:70:VAL:HG13	1:D:126:VAL:HG12	1.77	0.65
1:A:3:LYS:HG2	1:A:78:VAL:HB	1.78	0.64
1:A:358:LYS:NZ	1:A:361:GLU:OE1	2.23	0.64
1:B:20:ARG:HH11	1:B:354:THR:HG21	1.62	0.64
1:C:339:ILE:HG13	1:C:407:CYS:HB3	1.79	0.64
1:B:87:GLY:HA3	1:B:140:HIS:HB2	1.80	0.63
1:A:117:ASP:OD1	1:A:120:ARG:NH2	2.32	0.63
1:C:20:ARG:HH11	1:C:354:THR:HG21	1.64	0.63
1:B:120:ARG:NH1	1:B:334:LEU:HG	2.14	0.63
1:B:20:ARG:NH1	1:B:354:THR:HG21	2.14	0.62
1:B:246:GLU:HG2	1:B:249:LYS:HE3	1.81	0.62
1:C:246:GLU:HG2	1:C:249:LYS:HE3	1.82	0.62
1:D:358:LYS:HD2	1:D:359:PRO:HD2	1.80	0.61
1:C:20:ARG:NH1	1:C:354:THR:HG21	2.16	0.61
1:C:101:GLY:C	1:C:102:PHE:HD1	2.09	0.61
1:D:246:GLU:HG2	1:D:249:LYS:HE3	1.81	0.61
1:D:215:GLU:HB2	1:D:397:GLN:O	2.01	0.61
1:B:101:GLY:C	1:B:102:PHE:HD1	2.10	0.60
1:B:344:CYS:HB2	1:B:348:GLN:HG3	1.84	0.60
1:D:344:CYS:HB2	1:D:348:GLN:HG3	1.83	0.60
1:A:101:GLY:C	1:A:102:PHE:HD1	2.10	0.60
1:D:101:GLY:C	1:D:102:PHE:HD1	2.10	0.60
1:B:281:ASN:HA	1:B:284:GLU:HG3	1.83	0.59
1:C:220:ILE:HG12	1:C:355:PRO:HA	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:ASN:HA	1:A:284:GLU:HG3	1.84	0.59
1:C:281:ASN:HA	1:C:284:GLU:HG3	1.83	0.59
1:A:246:GLU:HG2	1:A:249:LYS:HE3	1.84	0.59
1:B:309:VAL:HG12	1:B:313:MET:HG2	1.85	0.58
1:D:281:ASN:HA	1:D:284:GLU:HG3	1.85	0.58
1:D:150:GLU:O	1:D:384:ARG:NH2	2.36	0.58
1:A:373:ILE:O	1:A:377:THR:HG23	2.04	0.58
1:A:309:VAL:HG12	1:A:313:MET:HG2	1.85	0.58
1:D:131:ARG:NH2	2:D:501:SO4:O1	2.35	0.58
1:B:387:VAL:HG21	1:B:405:LEU:HB2	1.85	0.57
1:D:321:GLU:CB	1:D:329:ARG:HD2	2.34	0.57
1:D:276:LYS:HE2	1:D:287:LEU:HD12	1.87	0.57
1:B:373:ILE:O	1:B:377:THR:HG23	2.04	0.57
1:C:373:ILE:O	1:C:377:THR:HG23	2.04	0.57
1:C:120:ARG:HD2	1:C:332:VAL:O	2.03	0.57
1:D:309:VAL:HG12	1:D:313:MET:HG2	1.87	0.57
1:C:309:VAL:HG12	1:C:313:MET:HG2	1.86	0.57
1:D:387:VAL:HG21	1:D:405:LEU:HB2	1.87	0.57
1:C:20:ARG:HD3	1:C:359:PRO:CB	2.29	0.57
1:C:280:GLY:O	1:C:282:GLY:N	2.38	0.56
1:A:344:CYS:HB2	1:A:348:GLN:HG3	1.86	0.56
1:C:276:LYS:HE2	1:C:287:LEU:HD12	1.87	0.56
1:C:132:MET:O	1:C:339:ILE:HG22	2.05	0.56
1:B:121:ASP:OD1	1:B:332:VAL:HG21	2.06	0.56
1:D:345:GLN:O	1:D:347:ASP:N	2.37	0.56
1:A:386:MET:HG2	1:A:416:PHE:CE1	2.40	0.56
1:D:121:ASP:OD1	1:D:332:VAL:HG21	2.06	0.55
1:B:276:LYS:HE2	1:B:287:LEU:HD12	1.88	0.55
1:C:23:VAL:HG13	1:C:48:ILE:HD13	1.89	0.55
1:C:111:MET:HE3	1:C:232:LEU:HD12	1.88	0.55
1:C:344:CYS:HB2	1:C:348:GLN:HG3	1.87	0.55
1:D:111:MET:HE2	1:D:235:ILE:HB	1.88	0.55
1:A:89:ARG:HD2	1:A:104:GLU:HG2	1.89	0.55
1:B:276:LYS:HE3	1:B:283:GLU:HB3	1.89	0.55
1:C:343:GLY:HA2	1:C:366:MET:HG3	1.89	0.55
1:A:139:CYS:SG	1:A:226:SER:HB2	2.47	0.55
1:B:386:MET:HG2	1:B:416:PHE:CE1	2.42	0.55
1:D:111:MET:HE3	1:D:232:LEU:HD12	1.88	0.55
1:A:132:MET:O	1:A:339:ILE:HG22	2.06	0.55
1:A:280:GLY:O	1:A:282:GLY:N	2.40	0.55
1:B:280:GLY:O	1:B:282:GLY:N	2.39	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:TYR:HD1	1:A:418:CYS:SG	2.28	0.55
1:B:37:TYR:HD1	1:B:418:CYS:SG	2.30	0.55
1:B:111:MET:HE3	1:B:232:LEU:HD12	1.89	0.55
1:C:386:MET:HG2	1:C:416:PHE:CE1	2.41	0.55
1:A:276:LYS:HE2	1:A:287:LEU:HD12	1.89	0.54
1:C:123:VAL:HG22	1:C:132:MET:SD	2.47	0.54
1:D:151:GLN:HB3	1:D:405:LEU:HB3	1.89	0.54
1:A:345:GLN:O	1:A:347:ASP:N	2.36	0.54
1:C:387:VAL:HG21	1:C:405:LEU:HB2	1.88	0.54
1:B:244:ASN:HD21	1:B:249:LYS:HZ2	1.56	0.54
1:C:37:TYR:HD1	1:C:418:CYS:SG	2.31	0.54
1:B:132:MET:O	1:B:339:ILE:HG22	2.07	0.54
1:D:386:MET:HG2	1:D:416:PHE:CE1	2.43	0.54
1:D:37:TYR:HD1	1:D:418:CYS:SG	2.30	0.54
1:D:280:GLY:O	1:D:282:GLY:N	2.40	0.54
1:B:89:ARG:HD2	1:B:104:GLU:HG2	1.89	0.53
1:A:111:MET:HE2	1:A:235:ILE:HB	1.90	0.53
1:A:343:GLY:HA2	1:A:366:MET:HG3	1.89	0.53
1:B:58:THR:OG1	1:B:61:ASN:OD1	2.27	0.53
1:C:276:LYS:HE3	1:C:283:GLU:HB3	1.89	0.53
1:D:120:ARG:NE	1:D:332:VAL:O	2.38	0.53
1:D:150:GLU:HG3	1:D:412:ALA:HB2	1.91	0.53
1:D:99:ASP:OD1	1:D:325:ALA:N	2.42	0.52
1:B:139:CYS:SG	1:B:226:SER:HB2	2.48	0.52
1:C:151:GLN:HB3	1:C:405:LEU:HB3	1.91	0.52
1:C:89:ARG:HD2	1:C:104:GLU:HG2	1.91	0.52
1:A:120:ARG:NH1	1:A:144:LEU:HA	2.25	0.52
1:A:344:CYS:HB2	1:A:348:GLN:HE21	1.75	0.52
1:B:343:GLY:HA2	1:B:366:MET:HG3	1.92	0.52
1:C:137:ASP:O	1:C:225:LYS:HE2	2.10	0.52
1:D:139:CYS:SG	1:D:226:SER:HB2	2.49	0.52
1:A:120:ARG:CZ	1:A:144:LEU:HA	2.39	0.52
1:C:210:LYS:O	1:C:222:ALA:N	2.42	0.52
1:A:244:ASN:HD21	1:A:249:LYS:HZ2	1.58	0.51
1:A:87:GLY:HA3	1:A:140:HIS:HB2	1.92	0.51
1:A:276:LYS:HE3	1:A:283:GLU:HB3	1.91	0.51
1:B:360:THR:HG21	1:C:238:GLN:HG2	1.91	0.51
1:D:344:CYS:HB2	1:D:348:GLN:CG	2.39	0.51
1:B:92:ALA:HB2	1:B:325:ALA:HB3	1.92	0.51
1:C:281:ASN:HA	1:C:284:GLU:CG	2.41	0.51
1:C:345:GLN:O	1:C:347:ASP:N	2.37	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:VAL:HG13	1:A:48:ILE:HD13	1.91	0.51
1:A:387:VAL:HG21	1:A:405:LEU:HB2	1.92	0.51
1:D:339:ILE:HG13	1:D:407:CYS:HB3	1.92	0.51
1:D:276:LYS:HE3	1:D:283:GLU:HB3	1.92	0.50
1:D:372:THR:O	1:D:376:GLU:HG3	2.10	0.50
1:A:151:GLN:HB3	1:A:405:LEU:HB3	1.93	0.50
1:B:255:PHE:CG	1:B:270:MET:HG3	2.46	0.50
1:B:345:GLN:O	1:B:347:ASP:N	2.36	0.50
1:C:121:ASP:CG	1:C:332:VAL:HG21	2.36	0.50
1:D:255:PHE:CG	1:D:270:MET:HG3	2.46	0.50
1:C:255:PHE:CG	1:C:270:MET:HG3	2.47	0.50
1:D:321:GLU:HB3	1:D:329:ARG:HD2	1.92	0.50
1:D:219:THR:O	1:D:355:PRO:O	2.29	0.50
1:A:58:THR:OG1	1:A:61:ASN:OD1	2.27	0.50
1:B:344:CYS:HB2	1:B:348:GLN:HE21	1.76	0.50
1:C:123:VAL:O	1:C:126:VAL:HG13	2.12	0.50
1:A:99:ASP:OD1	1:A:325:ALA:N	2.41	0.50
1:B:281:ASN:HA	1:B:284:GLU:CG	2.42	0.50
1:B:344:CYS:HB2	1:B:348:GLN:CG	2.41	0.50
1:C:281:ASN:HA	1:C:284:GLU:HB2	1.94	0.50
1:B:111:MET:HE2	1:B:235:ILE:HB	1.94	0.49
1:A:255:PHE:CG	1:A:270:MET:HG3	2.47	0.49
1:A:344:CYS:HB2	1:A:348:GLN:CG	2.42	0.49
1:C:382:SER:OG	1:C:385:GLU:HG3	2.12	0.49
1:A:382:SER:OG	1:A:385:GLU:HG3	2.12	0.49
1:A:123:VAL:HG22	1:A:132:MET:SD	2.52	0.49
1:A:244:ASN:HD21	1:A:249:LYS:NZ	2.11	0.49
1:B:244:ASN:HD21	1:B:249:LYS:NZ	2.10	0.49
1:D:89:ARG:HD2	1:D:104:GLU:HG2	1.93	0.49
1:D:102:PHE:O	1:D:326:GLY:HA2	2.13	0.49
1:C:244:ASN:HD21	1:C:249:LYS:HZ2	1.59	0.49
1:D:343:GLY:HA2	1:D:366:MET:HG3	1.94	0.49
1:A:111:MET:HE3	1:A:232:LEU:HD12	1.94	0.49
1:C:111:MET:HE2	1:C:235:ILE:HB	1.95	0.49
1:A:130:CYS:O	1:A:132:MET:N	2.46	0.49
1:A:222:ALA:HA	1:A:352:ASP:O	2.12	0.49
1:B:23:VAL:HG13	1:B:48:ILE:HD13	1.94	0.49
1:B:382:SER:OG	1:B:385:GLU:HG3	2.13	0.49
1:C:344:CYS:HB2	1:C:348:GLN:CG	2.42	0.49
1:B:123:VAL:HG22	1:B:132:MET:SD	2.52	0.49
1:C:344:CYS:HB2	1:C:348:GLN:HE21	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:12:TYR:OH	1:D:107:VAL:O	2.29	0.49
1:A:339:ILE:HD12	1:A:411:TYR:HB3	1.95	0.48
1:D:123:VAL:HG22	1:D:132:MET:SD	2.53	0.48
1:D:244:ASN:HD21	1:D:249:LYS:NZ	2.11	0.48
1:B:281:ASN:HA	1:B:284:GLU:HB2	1.95	0.48
1:C:90:LEU:HD11	1:C:250:ILE:HG21	1.95	0.48
1:D:63:ARG:NH1	1:D:122:LEU:HD11	2.28	0.48
1:D:321:GLU:HG3	1:D:324:TYR:CE2	2.48	0.48
1:B:151:GLN:HB3	1:B:405:LEU:HB3	1.95	0.48
1:A:281:ASN:HA	1:A:284:GLU:CG	2.43	0.48
1:B:339:ILE:HD12	1:B:411:TYR:HB3	1.95	0.48
1:C:89:ARG:HB3	1:C:102:PHE:HD2	1.79	0.48
1:C:139:CYS:SG	1:C:226:SER:HB2	2.52	0.48
1:D:281:ASN:HA	1:D:284:GLU:CG	2.43	0.48
1:D:132:MET:O	1:D:339:ILE:HG22	2.13	0.48
1:D:281:ASN:HA	1:D:284:GLU:HB2	1.94	0.48
1:A:236:LEU:HD23	1:A:236:LEU:HA	1.73	0.48
1:B:212:ILE:HB	1:B:221:HIS:CD2	2.48	0.48
1:B:89:ARG:HB3	1:B:102:PHE:HD2	1.79	0.48
1:A:150:GLU:O	1:A:384:ARG:NH2	2.47	0.48
1:C:115:THR:HG22	1:C:117:ASP:H	1.79	0.47
1:C:121:ASP:OD1	1:C:332:VAL:HG21	2.13	0.47
1:C:130:CYS:O	1:C:132:MET:N	2.48	0.47
1:C:87:GLY:HA3	1:C:140:HIS:HB2	1.96	0.47
1:C:385:GLU:O	1:C:389:ARG:HG3	2.13	0.47
1:D:385:GLU:O	1:D:389:ARG:HG3	2.14	0.47
1:A:385:GLU:O	1:A:389:ARG:HG3	2.14	0.47
1:B:240:THR:O	1:B:242:ASN:N	2.48	0.47
1:B:385:GLU:O	1:B:389:ARG:HG3	2.14	0.47
1:D:344:CYS:HB2	1:D:348:GLN:HE21	1.79	0.47
1:B:321:GLU:HG3	1:B:324:TYR:CE2	2.49	0.47
1:C:58:THR:OG1	1:C:61:ASN:OD1	2.31	0.47
1:D:339:ILE:HA	1:D:407:CYS:HB3	1.96	0.47
1:D:355:PRO:HD3	1:D:362:ALA:HA	1.96	0.47
1:A:120:ARG:HD3	1:A:144:LEU:HA	1.96	0.47
1:A:281:ASN:HA	1:A:284:GLU:HB2	1.96	0.47
1:B:130:CYS:O	1:B:132:MET:N	2.47	0.47
1:C:131:ARG:HH21	1:C:418:CYS:HA	1.80	0.47
1:C:339:ILE:HD12	1:C:411:TYR:HB3	1.97	0.47
1:D:236:LEU:HD23	1:D:236:LEU:HA	1.73	0.47
1:D:90:LEU:HD11	1:D:250:ILE:HG21	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:126:VAL:HG11	1:D:132:MET:HE3	1.95	0.47
1:C:236:LEU:HD23	1:C:236:LEU:HA	1.71	0.47
1:D:339:ILE:HD12	1:D:411:TYR:HB3	1.97	0.46
1:D:244:ASN:HD21	1:D:249:LYS:HZ2	1.63	0.46
1:B:238:GLN:CD	1:C:357:GLY:HA3	2.40	0.46
1:D:291:LEU:O	1:D:295:ALA:HB3	2.15	0.46
1:A:144:LEU:HD23	1:A:334:LEU:HD21	1.98	0.46
1:B:52:GLU:HG2	1:C:50:THR:O	2.16	0.46
1:A:212:ILE:HG13	1:A:221:HIS:CD2	2.51	0.46
1:A:240:THR:O	1:A:242:ASN:N	2.49	0.46
1:C:236:LEU:HB3	1:C:245:ILE:HD13	1.97	0.46
1:D:240:THR:O	1:D:242:ASN:N	2.49	0.46
1:D:291:LEU:HD23	1:D:291:LEU:HA	1.71	0.46
1:B:89:ARG:HH21	1:B:102:PHE:HB2	1.81	0.46
1:D:249:LYS:HZ1	1:D:292:GLY:HA3	1.81	0.46
1:D:384:ARG:O	1:D:388:THR:OG1	2.28	0.46
1:A:93:GLU:OE2	1:A:250:ILE:HB	2.16	0.45
1:A:122:LEU:HD23	1:A:122:LEU:HA	1.86	0.45
1:C:240:THR:O	1:C:242:ASN:N	2.49	0.45
1:A:5:ALA:HB3	1:A:44:ILE:HG12	1.98	0.45
1:D:58:THR:OG1	1:D:61:ASN:OD1	2.32	0.45
1:D:130:CYS:O	1:D:132:MET:N	2.49	0.45
1:C:123:VAL:HG12	1:C:335:PRO:HG2	1.97	0.45
1:C:244:ASN:HD21	1:C:249:LYS:NZ	2.14	0.45
1:A:291:LEU:O	1:A:295:ALA:HB3	2.17	0.45
1:C:281:ASN:ND2	1:C:284:GLU:OE2	2.50	0.45
1:D:123:VAL:O	1:D:126:VAL:HG13	2.17	0.45
1:B:116:ASP:O	1:B:120:ARG:HG3	2.15	0.45
1:B:150:GLU:O	1:B:384:ARG:NH2	2.49	0.45
1:D:89:ARG:HH21	1:D:102:PHE:HB2	1.81	0.45
1:A:89:ARG:HH21	1:A:102:PHE:HB2	1.82	0.45
1:C:16:LYS:HB2	1:C:16:LYS:HE2	1.55	0.45
1:C:150:GLU:O	1:C:384:ARG:NH2	2.49	0.45
1:C:228:PRO:HG2	1:C:231:THR:OG1	2.17	0.45
1:B:137:ASP:O	1:B:225:LYS:HE2	2.17	0.45
1:C:118:ASP:OD1	1:C:264:PRO:HB2	2.16	0.45
1:C:332:VAL:HA	1:C:333:PRO:HD2	1.78	0.45
1:D:50:THR:OG1	1:D:53:SER:HB3	2.17	0.45
1:B:126:VAL:HG11	1:B:132:MET:HE3	1.99	0.44
1:C:294:LEU:HD12	1:C:294:LEU:HA	1.70	0.44
1:A:228:PRO:HG2	1:A:231:THR:OG1	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:5:ALA:HB3	1:C:44:ILE:HG12	1.99	0.44
1:D:281:ASN:ND2	1:D:284:GLU:OE2	2.50	0.44
1:D:294:LEU:HD12	1:D:294:LEU:HA	1.71	0.44
1:B:294:LEU:HD12	1:B:294:LEU:HA	1.72	0.44
1:A:89:ARG:HB3	1:A:102:PHE:HD2	1.81	0.44
1:A:281:ASN:ND2	1:A:284:GLU:OE2	2.51	0.44
1:C:12:TYR:OH	1:C:107:VAL:O	2.31	0.44
1:D:131:ARG:HH21	1:D:418:CYS:HA	1.83	0.44
1:C:99:ASP:HB2	1:C:324:TYR:CG	2.52	0.44
1:C:20:ARG:NH1	2:C:501:SO4:O4	2.50	0.44
1:B:5:ALA:HB3	1:B:44:ILE:HG12	1.99	0.44
1:C:140:HIS:CE1	1:C:226:SER:OG	2.71	0.44
1:C:374:LEU:HD23	1:C:374:LEU:HA	1.75	0.44
1:A:12:TYR:OH	1:A:107:VAL:O	2.25	0.43
1:A:325:ALA:C	1:A:327:GLY:H	2.25	0.43
1:A:321:GLU:HG3	1:A:324:TYR:CE2	2.54	0.43
1:C:126:VAL:HG11	1:C:132:MET:HE3	2.00	0.43
1:C:372:THR:O	1:C:376:GLU:HG3	2.17	0.43
1:D:212:ILE:HB	1:D:221:HIS:HD2	1.84	0.43
1:A:372:THR:O	1:A:376:GLU:HG3	2.18	0.43
1:B:90:LEU:HD11	1:B:250:ILE:HG21	1.99	0.43
1:C:210:LYS:HB2	1:C:221:HIS:HB3	2.00	0.43
1:A:92:ALA:HB2	1:A:325:ALA:HB3	2.00	0.43
1:A:396:LYS:HG3	1:A:397:GLN:HE21	1.83	0.43
1:B:6:VAL:HG23	1:B:69:LEU:HD22	2.00	0.43
1:B:236:LEU:HB3	1:B:245:ILE:HD13	2.00	0.43
1:C:291:LEU:O	1:C:295:ALA:HB3	2.19	0.43
1:D:124:ASP:HA	1:D:335:PRO:HG2	2.00	0.43
1:B:131:ARG:HH21	1:B:418:CYS:HA	1.84	0.43
1:B:281:ASN:ND2	1:B:284:GLU:OE2	2.51	0.43
1:B:349:THR:O	1:B:401:GLN:HB3	2.19	0.43
1:C:291:LEU:HD23	1:C:291:LEU:HA	1.73	0.43
1:D:137:ASP:O	1:D:225:LYS:HE2	2.19	0.43
1:A:126:VAL:HG11	1:A:132:MET:HE3	2.01	0.43
1:B:30:TYR:OH	1:B:41:GLU:HB3	2.18	0.43
1:B:372:THR:O	1:B:376:GLU:HG3	2.18	0.43
1:D:49:ASP:CG	1:D:50:THR:N	2.77	0.43
1:A:19:LEU:HD23	1:A:223:LYS:HD3	2.01	0.43
1:A:30:TYR:OH	1:A:41:GLU:HB3	2.19	0.43
1:B:150:GLU:HG3	1:B:412:ALA:HB2	2.01	0.43
1:C:116:ASP:O	1:C:120:ARG:HG3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:333:PRO:O	1:D:335:PRO:HD3	2.19	0.43
1:D:354:THR:HA	1:D:355:PRO:HD3	1.87	0.43
1:A:236:LEU:HB3	1:A:245:ILE:HD13	2.00	0.42
1:C:349:THR:O	1:C:401:GLN:HB3	2.18	0.42
1:D:122:LEU:HD23	1:D:122:LEU:HA	1.88	0.42
1:D:332:VAL:HA	1:D:333:PRO:HD2	1.66	0.42
1:D:377:THR:H	1:D:377:THR:HG23	1.60	0.42
1:C:260:ASP:O	1:C:267:LYS:HE2	2.19	0.42
1:D:382:SER:OG	1:D:385:GLU:HG3	2.19	0.42
1:A:349:THR:O	1:A:401:GLN:HB3	2.19	0.42
1:D:349:THR:O	1:D:401:GLN:HB3	2.19	0.42
1:C:93:GLU:OE2	1:C:250:ILE:HB	2.19	0.42
1:D:5:ALA:HB3	1:D:44:ILE:HG12	2.01	0.42
1:A:260:ASP:O	1:A:267:LYS:HE2	2.20	0.42
1:B:228:PRO:HG2	1:B:231:THR:OG1	2.19	0.42
1:C:70:VAL:HG13	1:C:126:VAL:CG1	2.46	0.42
1:C:99:ASP:OD1	1:C:325:ALA:N	2.53	0.42
1:D:6:VAL:HG23	1:D:69:LEU:HD22	2.01	0.42
1:D:228:PRO:HG2	1:D:231:THR:OG1	2.20	0.42
1:B:122:LEU:HD23	1:B:122:LEU:HA	1.91	0.42
1:A:90:LEU:HD11	1:A:250:ILE:HG21	2.01	0.42
1:C:145:ILE:HG21	1:C:406:TYR:CB	2.46	0.42
1:C:220:ILE:HG23	1:C:354:THR:H	1.85	0.42
1:A:123:VAL:O	1:A:126:VAL:HG13	2.20	0.41
1:A:137:ASP:O	1:A:225:LYS:HE2	2.20	0.41
1:D:260:ASP:O	1:D:267:LYS:HE2	2.20	0.41
1:A:249:LYS:HZ1	1:A:292:GLY:HA3	1.85	0.41
1:B:236:LEU:HD23	1:B:236:LEU:HA	1.77	0.41
1:A:211:GLU:O	1:A:212:ILE:C	2.63	0.41
1:D:93:GLU:OE2	1:D:250:ILE:HB	2.20	0.41
1:B:93:GLU:OE2	1:B:250:ILE:HB	2.19	0.41
1:C:30:TYR:OH	1:C:41:GLU:HB3	2.21	0.41
1:B:99:ASP:OD1	1:B:325:ALA:N	2.45	0.41
1:B:288:MET:HE2	1:B:288:MET:HB3	1.83	0.41
1:C:249:LYS:HZ1	1:C:292:GLY:HA3	1.85	0.41
1:D:30:TYR:OH	1:D:41:GLU:HB3	2.20	0.41
1:D:396:LYS:HG3	1:D:397:GLN:HE21	1.85	0.41
1:D:130:CYS:O	1:D:337:SER:OG	2.39	0.41
1:A:332:VAL:HA	1:A:333:PRO:HD2	1.80	0.41
1:B:239:GLN:HA	1:C:358:LYS:CE	2.51	0.41
1:C:6:VAL:HG23	1:C:69:LEU:HD22	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:233:ILE:O	1:C:237:LYS:HG3	2.21	0.41
1:A:221:HIS:O	1:A:354:THR:HG22	2.20	0.41
1:A:291:LEU:HD23	1:A:291:LEU:HA	1.73	0.41
1:B:92:ALA:HA	1:B:325:ALA:HB2	2.02	0.41
1:B:124:ASP:HA	1:B:335:PRO:HG2	2.02	0.41
1:B:260:ASP:O	1:B:267:LYS:HE2	2.21	0.41
1:B:18:GLU:O	1:B:223:LYS:HE2	2.20	0.41
1:B:121:ASP:CG	1:B:332:VAL:HG21	2.45	0.41
1:B:123:VAL:O	1:B:126:VAL:HG13	2.21	0.41
1:C:89:ARG:HH21	1:C:102:PHE:HB2	1.86	0.41
1:C:150:GLU:HG3	1:C:412:ALA:HB2	2.01	0.41
1:A:145:ILE:HG12	1:A:406:TYR:CD2	2.56	0.40
1:A:224:ASP:C	1:A:226:SER:H	2.29	0.40
1:D:236:LEU:HB3	1:D:245:ILE:HD13	2.03	0.40
1:B:4:LYS:O	1:B:79:LEU:HA	2.21	0.40
1:B:373:ILE:HD11	1:B:390:ALA:HA	2.03	0.40
1:A:150:GLU:HG3	1:A:412:ALA:HB2	2.04	0.40
1:D:10:ILE:HG12	1:D:57:PRO:CG	2.50	0.40
1:D:70:VAL:HG13	1:D:126:VAL:CG1	2.49	0.40
1:A:294:LEU:HA	1:A:294:LEU:HD12	1.68	0.40
1:A:373:ILE:HD11	1:A:390:ALA:HA	2.03	0.40
1:C:104:GLU:HB3	1:C:140:HIS:O	2.21	0.40
1:C:321:GLU:HG3	1:C:324:TYR:CE2	2.57	0.40
1:D:288:MET:HE2	1:D:288:MET:HB3	1.84	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	358/426 (84%)	324 (90%)	28 (8%)	6 (2%)	7 34

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	351/426 (82%)	318 (91%)	27 (8%)	6 (2%)	7	34
1	C	355/426 (83%)	320 (90%)	28 (8%)	7 (2%)	6	32
1	D	352/426 (83%)	319 (91%)	25 (7%)	8 (2%)	5	30
All	All	1416/1704 (83%)	1281 (90%)	108 (8%)	27 (2%)	6	32

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	131	ARG
1	A	212	ILE
1	A	327	GLY
1	A	346	THR
1	A	356	ALA
1	B	131	ARG
1	B	327	GLY
1	B	346	THR
1	B	356	ALA
1	C	131	ARG
1	C	346	THR
1	C	356	ALA
1	D	50	THR
1	D	131	ARG
1	D	346	THR
1	D	360	THR
1	A	241	GLY
1	B	212	ILE
1	B	241	GLY
1	C	210	LYS
1	C	241	GLY
1	D	241	GLY
1	D	356	ALA
1	C	146	ASP
1	C	208	GLU
1	D	146	ASP
1	D	339	ILE

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	300/357 (84%)	289 (96%)	11 (4%)	30	56
1	B	297/357 (83%)	285 (96%)	12 (4%)	28	54
1	C	301/357 (84%)	289 (96%)	12 (4%)	28	54
1	D	299/357 (84%)	288 (96%)	11 (4%)	30	56
All	All	1197/1428 (84%)	1151 (96%)	46 (4%)	29	56

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	VAL
1	A	123	VAL
1	A	141	SER
1	A	145	ILE
1	A	151	GLN
1	A	209	THR
1	A	218	GLU
1	A	236	LEU
1	A	283	GLU
1	A	300	GLU
1	A	381	ILE
1	B	16	LYS
1	B	26	VAL
1	B	123	VAL
1	B	151	GLN
1	B	209	THR
1	B	211	GLU
1	B	221	HIS
1	B	236	LEU
1	B	283	GLU
1	B	293	LYS
1	B	300	GLU
1	B	381	ILE
1	C	16	LYS
1	C	26	VAL
1	C	123	VAL
1	C	145	ILE
1	C	151	GLN
1	C	206	GLU
1	C	208	GLU

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Mol	Chain	Res	Type
1	C	209	THR
1	C	236	LEU
1	C	283	GLU
1	C	300	GLU
1	C	381	ILE
1	D	26	VAL
1	D	123	VAL
1	D	145	ILE
1	D	151	GLN
1	D	209	THR
1	D	210	LYS
1	D	218	GLU
1	D	236	LEU
1	D	283	GLU
1	D	300	GLU
1	D	381	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	221	HIS
1	A	244	ASN
1	A	281	ASN
1	A	314	GLN
1	A	348	GLN
1	A	397	GLN
1	A	402	GLN
1	B	244	ASN
1	B	281	ASN
1	B	314	GLN
1	B	348	GLN
1	B	402	GLN
1	C	140	HIS
1	C	221	HIS
1	C	244	ASN
1	C	281	ASN
1	C	314	GLN
1	C	348	GLN
1	D	56	GLN
1	D	244	ASN
1	D	281	ASN
1	D	314	GLN

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Mol	Chain	Res	Type
1	D	348	GLN
1	D	397	GLN

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](#)

8 ligands are modelled in this entry.

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	SO4	B	501	-	4,4,4	0.24	0	6,6,6	0.07	0
2	SO4	C	502	-	4,4,4	0.24	0	6,6,6	0.07	0
2	SO4	D	501	-	4,4,4	0.24	0	6,6,6	0.09	0
2	SO4	C	501	-	4,4,4	0.24	0	6,6,6	0.06	0
2	SO4	A	503	-	4,4,4	0.24	0	6,6,6	0.07	0
2	SO4	A	501	-	4,4,4	0.22	0	6,6,6	0.15	0
2	SO4	C	503	-	4,4,4	0.24	0	6,6,6	0.08	0
2	SO4	A	502	-	4,4,4	0.23	0	6,6,6	0.09	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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	501	SO4	1	0
2	C	501	SO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	362/426 (84%)	0.25	15 (4%) 41 23	43, 86, 129, 185	0
1	B	357/426 (83%)	0.26	10 (2%) 55 32	50, 92, 144, 216	0
1	C	361/426 (84%)	0.26	11 (3%) 52 31	41, 84, 136, 182	0
1	D	358/426 (84%)	0.36	10 (2%) 55 32	49, 102, 150, 196	0
All	All	1438/1704 (84%)	0.28	46 (3%) 50 29	41, 90, 141, 216	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	213	GLU	4.2
1	D	151	GLN	3.7
1	B	213	GLU	3.6
1	A	214	LEU	3.5
1	D	293	LYS	3.5
1	A	144	LEU	3.4
1	A	212	ILE	3.4
1	A	213	GLU	3.4
1	C	151	GLN	3.3
1	A	420	GLU	3.3
1	A	143	GLY	3.2
1	A	151	GLN	3.2
1	C	337	SER	3.2
1	B	214	LEU	3.1
1	B	420	GLU	3.1
1	B	148	ALA	2.9
1	A	145	ILE	2.9
1	D	49	ASP	2.9
1	B	151	GLN	2.8
1	A	330	GLY	2.8
1	C	295	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	358	LYS	2.7
1	C	148	ALA	2.7
1	D	420	GLU	2.6
1	D	408	HIS	2.6
1	B	146	ASP	2.6
1	C	294	LEU	2.5
1	C	296	SER	2.4
1	A	215	GLU	2.4
1	C	293	LYS	2.4
1	D	337	SER	2.3
1	D	339	ILE	2.3
1	B	272	VAL	2.3
1	D	51	ASP	2.3
1	A	150	GLU	2.2
1	D	148	ALA	2.2
1	C	378	ASP	2.2
1	C	145	ILE	2.2
1	A	148	ALA	2.2
1	B	330	GLY	2.2
1	B	277	LEU	2.1
1	C	420	GLU	2.1
1	C	206	GLU	2.1
1	A	302	LYS	2.0
1	A	334	LEU	2.0
1	B	337	SER	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 ⓘ

There are no oligosaccharides 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. 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($\text{\AA}^2$)	Q<0.9
2	SO4	A	501	5/5	0.64	0.19	126,126,126,126	0
2	SO4	C	503	5/5	0.65	0.19	153,153,153,153	0
2	SO4	D	501	5/5	0.72	0.14	141,141,141,141	0
2	SO4	A	502	5/5	0.73	0.13	143,143,143,143	0
2	SO4	A	503	5/5	0.80	0.14	137,137,137,137	0
2	SO4	C	502	5/5	0.82	0.10	118,118,118,118	0
2	SO4	B	501	5/5	0.83	0.17	117,117,117,117	0
2	SO4	C	501	5/5	0.89	0.13	116,116,116,116	0

6.5 Other polymers [i](#)

There are no such residues in this entry.