



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 12:31 AM UTC

PDB ID : 1RFZ / pdb_00001rfz
Title : Structure of Protein of Unknown Function from Bacillus stearothermophilus
Authors : Kim, Y.; Wu, R.; Cuff, M.E.; Quartey, P.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2003-11-10
Resolution : 2.80 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

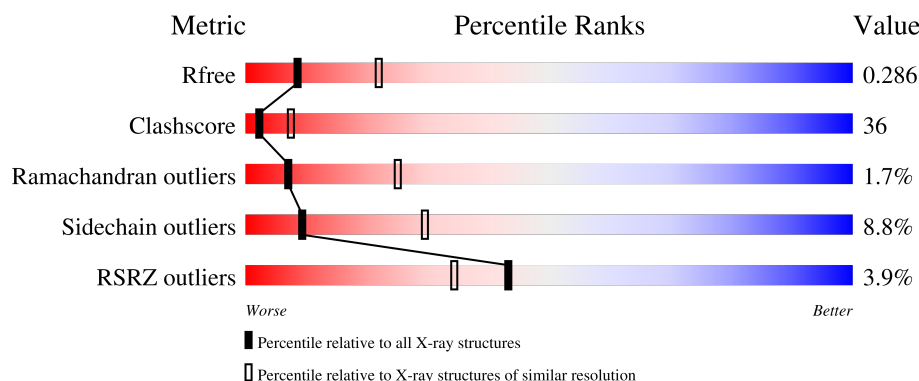
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	168	36% 54% 8% .
1	B	168	2% 47% 42% 8% ..
1	C	168	5% 48% 40% 7% ..
1	D	168	7% 28% 54% 8% . 8%

2 Entry composition [i](#)

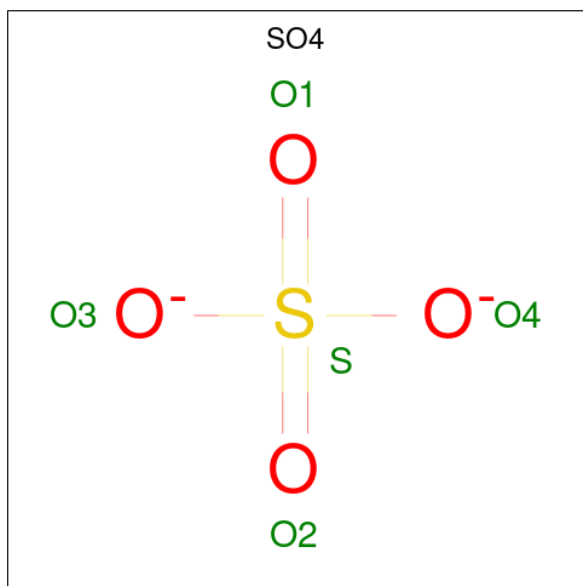
There are 3 unique types of molecules in this entry. The entry contains 5264 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Hypothetical protein APC35681.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	164	Total	C	N	O	S	Se	0	0	0
			1295	812	224	254	2	3			
1	B	164	Total	C	N	O	S	Se	0	2	0
			1317	824	232	256	2	3			
1	C	161	Total	C	N	O	S	Se	0	0	0
			1273	800	220	249	2	2			
1	D	154	Total	C	N	O	S	Se	0	1	0
			1227	770	213	241	2	1			

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



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

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

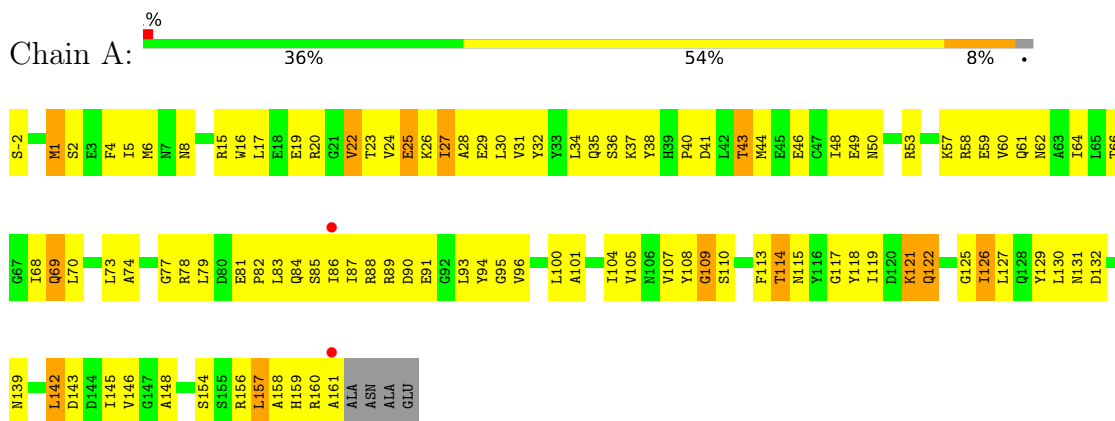
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	28	Total	O	0	0
			28	28		
3	B	44	Total	O	0	0
			44	44		
3	C	31	Total	O	0	0
			31	31		
3	D	14	Total	O	0	0
			14	14		

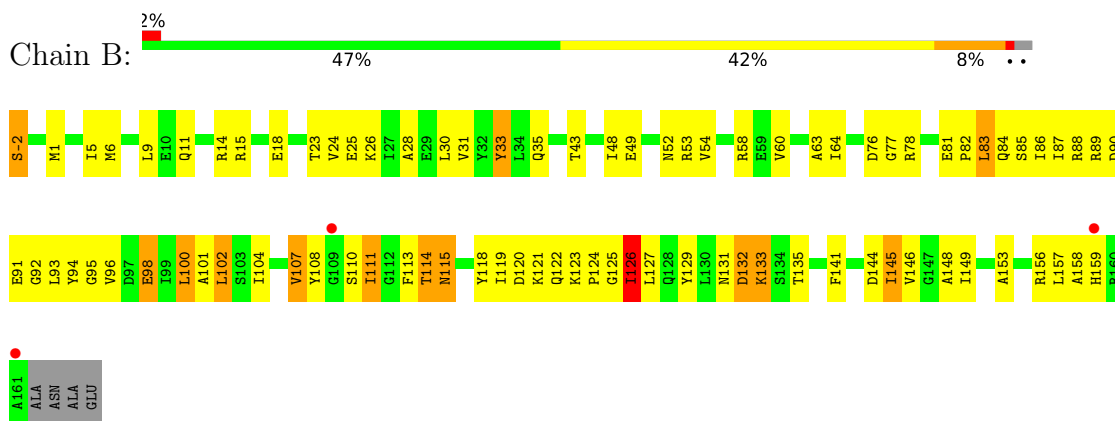
3 Residue-property plots

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.

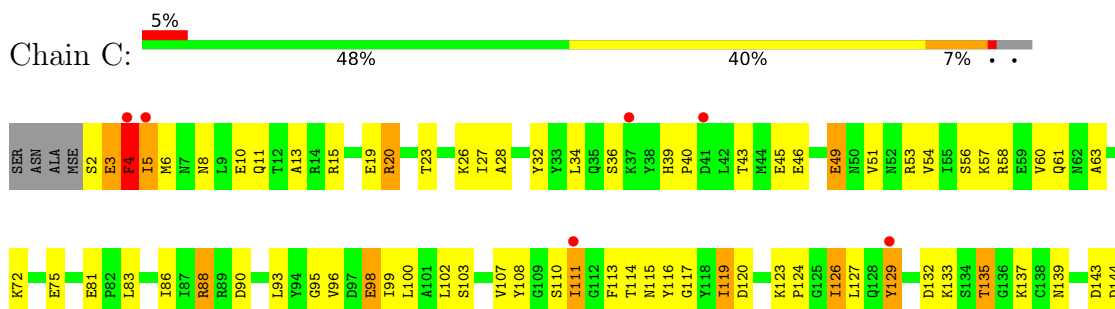
• Molecule 1: Hypothetical protein APC35681

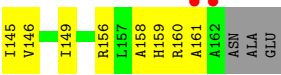


• Molecule 1: Hypothetical protein APC35681

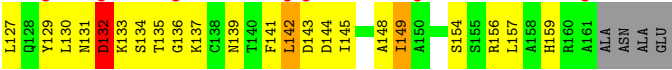
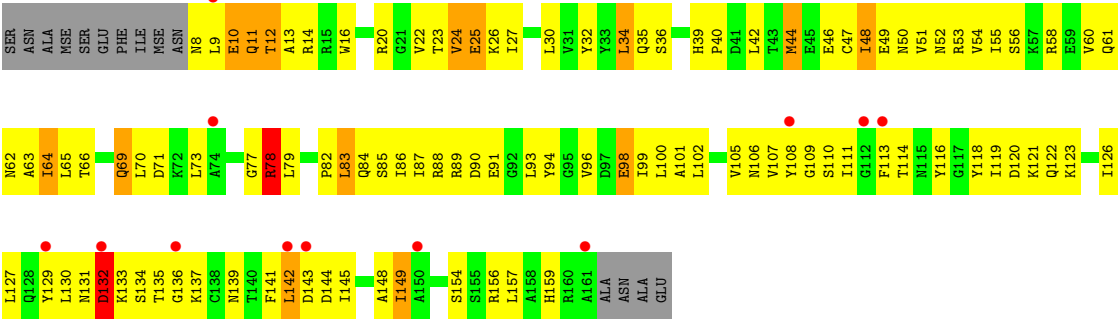


• Molecule 1: Hypothetical protein APC35681





● Molecule 1: Hypothetical protein APC35681



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	80.82Å 80.82Å 236.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.77 – 2.80 47.77 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.5 (47.77-2.80) 97.5 (47.77-2.80)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.26 (at 2.69Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.221 , 0.277 0.229 , 0.286	Depositor DCC
R_{free} test set	3568 reflections (8.64%)	wwPDB-VP
Wilson B-factor (Å ²)	56.1	Xtriage
Anisotropy	0.302	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 59.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5264	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.24% 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	0.48	0/1309	1.03	6/1763 (0.3%)
1	B	0.59	0/1331	0.98	6/1791 (0.3%)
1	C	0.59	0/1288	1.08	7/1737 (0.4%)
1	D	0.44	0/1242	1.04	8/1677 (0.5%)
All	All	0.53	0/5170	1.03	27/6968 (0.4%)

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	4	PHE	N-CA-C	-11.32	97.97	112.23
1	D	131	ASN	N-CA-C	-9.71	102.67	112.97
1	A	109	GLY	N-CA-C	9.18	121.40	112.04
1	C	3	GLU	N-CA-C	-9.13	103.29	112.97
1	D	50	ASN	N-CA-C	-8.78	101.65	111.14
1	C	135	THR	N-CA-C	-7.34	103.46	113.30
1	A	61	GLN	N-CA-C	-7.12	103.45	111.14
1	B	92	GLY	N-CA-C	6.99	122.62	114.16
1	A	142	LEU	N-CA-C	6.61	119.33	111.33
1	D	78	ARG	N-CA-C	-6.43	105.39	112.72
1	A	27	ILE	N-CA-C	-6.15	104.75	110.53
1	D	109	GLY	N-CA-C	6.12	119.74	112.33
1	B	31	VAL	N-CA-C	-6.10	104.39	110.62
1	B	146	VAL	N-CA-C	-5.89	104.99	110.53
1	B	153	ALA	N-CA-C	-5.56	105.12	111.07
1	A	22	VAL	N-CA-C	5.51	116.51	108.58
1	B	5	ILE	N-CA-C	5.50	115.70	110.53
1	C	45	GLU	N-CA-C	-5.48	105.20	111.07
1	D	77	GLY	N-CA-C	5.37	125.90	113.18
1	D	44	MSE	N-CA-C	-5.29	106.08	112.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	48	ILE	N-CA-C	5.29	116.64	110.62
1	C	98	GLU	N-CA-C	-5.19	105.62	111.28
1	C	39	HIS	CA-C-N	5.16	126.28	119.84
1	C	39	HIS	C-N-CA	5.16	126.28	119.84
1	D	16	TRP	N-CA-C	-5.10	105.80	111.36
1	B	126	ILE	CB-CA-C	-5.07	105.29	112.14
1	A	1	MSE	N-CA-C	-5.07	105.94	111.82

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	1295	0	1296	124	0
1	B	1317	0	1321	88	0
1	C	1273	0	1273	88	0
1	D	1227	0	1227	135	0
2	A	15	0	0	1	0
2	B	5	0	0	0	0
2	C	10	0	0	0	0
2	D	5	0	0	1	0
3	A	28	0	0	5	0
3	B	44	0	0	6	0
3	C	31	0	0	8	0
3	D	14	0	0	3	0
All	All	5264	0	5117	374	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 36.

All (374) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:MSE:HE1	1:D:86:ILE:HG12	1.25	1.17

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:86:ILE:HG13	1:C:6:MSE:HE3	1.26	1.14
1:A:70:LEU:HD21	1:D:70:LEU:HD21	1.45	0.97
1:C:100:LEU:O	1:C:103:SER:HB3	1.66	0.96
1:A:23:THR:HG22	1:A:25:GLU:H	1.32	0.94
1:B:23:THR:HG22	1:B:25:GLU:H	1.32	0.93
1:A:94:TYR:HA	1:D:62:ASN:HD21	1.34	0.92
1:D:14:ARG:HB2	3:D:417:HOH:O	1.71	0.90
1:D:85:SER:O	1:D:89:ARG:HG3	1.74	0.88
1:A:69:GLN:HE21	1:A:69:GLN:HA	1.39	0.87
1:B:86:ILE:CG1	1:C:6:MSE:HE3	2.03	0.87
1:A:32:TYR:O	1:A:36:SER:HB2	1.76	0.86
1:D:142:LEU:HA	1:D:145:ILE:HG12	1.60	0.84
1:A:5:ILE:HG22	1:A:6:MSE:HE2	1.60	0.83
1:A:-2:SER:HA	1:D:89:ARG:HH22	1.41	0.83
1:A:6:MSE:HE1	1:D:86:ILE:CG1	2.06	0.83
1:C:98:GLU:HB3	1:D:110:SER:HB2	1.59	0.81
1:A:6:MSE:CE	1:D:86:ILE:HG12	2.10	0.81
1:D:96:VAL:HA	1:D:99:ILE:HD12	1.62	0.80
1:D:42:LEU:HD11	1:D:46:GLU:HG2	1.63	0.80
1:C:132:ASP:O	1:C:133:LYS:HB2	1.81	0.79
1:D:26:LYS:HE3	1:D:126:ILE:HD12	1.66	0.77
1:D:142:LEU:HA	1:D:145:ILE:CG1	2.15	0.77
1:A:25:GLU:O	1:A:29:GLU:HG3	1.86	0.76
1:A:23:THR:HG22	1:A:25:GLU:N	2.00	0.75
1:C:43:THR:OG1	1:C:46:GLU:HG2	1.86	0.75
1:B:86:ILE:HG13	1:C:6:MSE:CE	2.14	0.75
1:D:110:SER:HA	1:D:113:PHE:CD2	2.20	0.75
1:C:111:ILE:HD12	1:C:111:ILE:H	1.52	0.74
1:D:130:LEU:O	1:D:139:ASN:HB2	1.87	0.74
1:A:83:LEU:HD22	1:A:87:ILE:HD11	1.69	0.74
1:B:111:ILE:HD11	3:B:435:HOH:O	1.86	0.74
1:B:23:THR:HG22	1:B:25:GLU:N	2.01	0.74
1:C:34:LEU:HD21	1:C:119:ILE:HG23	1.69	0.73
1:D:110:SER:HA	1:D:113:PHE:CE2	2.23	0.73
1:B:91:GLU:HG2	1:C:3:GLU:O	1.89	0.73
1:B:132:ASP:O	1:B:133:LYS:HD3	1.89	0.73
1:A:108:TYR:CZ	1:A:156:ARG:HD2	2.23	0.72
1:A:93:LEU:HD13	1:D:58:ARG:CZ	2.20	0.72
1:A:101:ALA:O	1:A:105:VAL:HG23	1.90	0.70
1:A:53:ARG:HG2	1:A:156:ARG:CZ	2.21	0.69
1:A:85:SER:O	1:A:89:ARG:HG2	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:115:ASN:O	1:C:119:ILE:HG13	1.91	0.69
1:A:53:ARG:HG2	1:A:156:ARG:NE	2.08	0.69
1:D:139:ASN:HA	1:D:142:LEU:HD21	1.73	0.68
1:B:120:ASP:O	1:B:123:LYS:HE2	1.92	0.68
1:A:96:VAL:HG13	1:D:100:LEU:HD21	1.76	0.68
1:D:14:ARG:HH11	1:D:24:VAL:HG21	1.59	0.67
1:C:4:PHE:CD1	1:C:4:PHE:C	2.72	0.67
1:A:40:PRO:HG2	1:A:41:ASP:H	1.58	0.67
1:B:91:GLU:OE2	1:C:6:MSE:HB2	1.95	0.67
1:A:81:GLU:OE2	1:A:82:PRO:HA	1.94	0.67
3:B:423:HOH:O	1:D:111:ILE:HD13	1.95	0.67
1:B:26:LYS:HG3	3:B:427:HOH:O	1.95	0.66
1:A:122:GLN:HE21	1:A:122:GLN:HA	1.60	0.66
1:B:-2:SER:HB2	1:B:1:MSE:HB3	1.77	0.66
1:C:111:ILE:HD12	1:C:111:ILE:N	2.11	0.66
1:D:130:LEU:HD13	1:D:142:LEU:HD12	1.77	0.66
1:B:102:LEU:HG	1:B:113:PHE:HE1	1.60	0.65
1:B:133:LYS:HG2	1:B:133:LYS:O	1.95	0.65
1:D:42:LEU:HD11	1:D:46:GLU:CG	2.26	0.65
1:D:132:ASP:C	1:D:134:SER:H	2.04	0.65
1:D:9:LEU:O	1:D:10:GLU:C	2.39	0.65
1:D:44:MSE:O	1:D:48:ILE:HG13	1.95	0.65
1:D:26:LYS:CE	1:D:126:ILE:HD12	2.26	0.65
1:C:156:ARG:HH12	1:C:160:ARG:HH11	1.43	0.65
1:D:82:PRO:O	1:D:86:ILE:HG13	1.96	0.64
1:D:69:GLN:HA	1:D:69:GLN:HE21	1.62	0.64
1:B:89[A]:ARG:HB3	1:C:3:GLU:HG2	1.78	0.64
1:A:113:PHE:CD2	1:B:102:LEU:HD13	2.33	0.64
1:A:83:LEU:HD22	1:A:87:ILE:CD1	2.27	0.64
1:A:60:VAL:O	1:A:64:ILE:HG13	1.97	0.64
1:C:49:GLU:HA	3:C:416:HOH:O	1.97	0.63
1:A:158:ALA:O	1:C:4:PHE:CE2	2.52	0.63
1:D:71:ASP:OD2	1:D:142:LEU:HD23	1.98	0.63
1:D:156:ARG:NE	2:D:404:SO4:O4	2.29	0.63
1:C:114:THR:CG2	1:D:121:LYS:HB2	2.28	0.63
1:A:1:MSE:O	1:A:5:ILE:HG13	2.00	0.62
1:C:110:SER:HA	1:C:113:PHE:CE2	2.34	0.62
1:A:66:THR:HA	1:D:83:LEU:HD11	1.82	0.62
1:A:78:ARG:NH1	1:D:78:ARG:HG3	2.15	0.62
1:A:69:GLN:O	1:A:73:LEU:HD12	1.99	0.62
1:D:108:TYR:CE2	1:D:156:ARG:HD2	2.35	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:134:SER:C	1:D:136:GLY:H	2.08	0.61
1:C:119:ILE:HD12	1:C:120:ASP:N	2.14	0.61
1:C:90:ASP:OD1	1:C:90:ASP:O	2.19	0.61
1:C:113:PHE:CZ	1:D:102:LEU:HD11	2.35	0.60
1:A:110:SER:OG	1:B:95:GLY:HA3	2.01	0.60
1:C:75:GLU:OE1	1:C:137:LYS:HD3	2.00	0.60
1:A:104:ILE:O	1:A:107:VAL:HB	2.02	0.60
1:B:88[B]:ARG:HD2	3:B:404:HOH:O	2.00	0.60
1:C:95:GLY:HA2	3:C:431:HOH:O	2.02	0.60
1:B:81:GLU:OE2	1:B:82:PRO:HA	2.01	0.60
1:D:10:GLU:HG2	1:D:61:GLN:HE22	1.67	0.60
1:B:33:TYR:C	1:B:33:TYR:CD2	2.80	0.59
1:A:83:LEU:HD11	1:D:66:THR:HA	1.84	0.59
1:D:118:TYR:CE1	1:D:122:GLN:HG3	2.37	0.59
1:A:74:ALA:O	1:A:88:ARG:HD2	2.02	0.59
1:B:89[B]:ARG:HB3	1:B:89[B]:ARG:HH11	1.67	0.59
1:A:44:MSE:O	1:A:48:ILE:HG13	2.01	0.59
1:B:102:LEU:HG	1:B:113:PHE:CE1	2.37	0.59
1:C:129:TYR:CD1	1:C:129:TYR:C	2.80	0.59
1:D:10:GLU:HG2	1:D:61:GLN:NE2	2.18	0.59
1:D:135:THR:C	1:D:137:LYS:H	2.11	0.59
1:C:20:ARG:HH22	1:C:75:GLU:CD	2.11	0.59
1:B:77:GLY:HA2	1:B:84:GLN:NE2	2.18	0.59
1:A:78:ARG:HD2	1:D:78:ARG:HD3	1.85	0.58
1:D:71:ASP:OD2	1:D:141:PHE:HB2	2.03	0.58
1:A:114:THR:HG23	3:A:408:HOH:O	2.03	0.58
1:D:79:LEU:HB2	1:D:84:GLN:OE1	2.04	0.58
1:B:82:PRO:HB2	1:C:6:MSE:HE1	1.85	0.57
1:D:130:LEU:C	1:D:132:ASP:H	2.11	0.57
1:D:23:THR:HG22	1:D:26:LYS:NZ	2.19	0.57
1:D:23:THR:C	1:D:25:GLU:H	2.11	0.57
1:A:43:THR:O	1:A:46:GLU:HB3	2.04	0.57
1:C:111:ILE:H	1:C:111:ILE:CD1	2.16	0.57
1:B:89[A]:ARG:CB	1:C:3:GLU:HG2	2.35	0.57
1:A:142:LEU:O	1:A:146:VAL:HG23	2.05	0.56
1:D:11:GLN:O	1:D:12:THR:C	2.48	0.56
1:D:32:TYR:O	1:D:36:SER:HB2	2.04	0.56
1:D:49[B]:GLU:O	1:D:53:ARG:HG2	2.06	0.56
1:B:43:THR:HA	3:B:439:HOH:O	2.05	0.56
1:D:129:TYR:CE1	1:D:135:THR:HG21	2.40	0.56
1:C:102:LEU:HD11	1:D:113:PHE:CD2	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:156:ARG:HH12	1:C:160:ARG:NH1	2.04	0.56
1:A:95:GLY:H	1:D:62:ASN:ND2	2.03	0.56
1:C:19:GLU:HG2	3:C:435:HOH:O	2.06	0.56
1:B:110:SER:HA	1:B:113:PHE:CD2	2.41	0.56
1:A:35:GLN:HG3	1:A:157:LEU:HD23	1.87	0.56
1:B:93:LEU:HD21	1:C:4:PHE:CE2	2.41	0.55
1:B:35:GLN:HG3	1:B:157:LEU:HD23	1.89	0.55
1:C:23:THR:OG1	1:C:26:LYS:HG2	2.07	0.55
1:A:142:LEU:HD12	1:A:145:ILE:HD11	1.87	0.55
1:A:50:ASN:OD1	1:A:156:ARG:NH1	2.39	0.55
1:D:110:SER:HA	1:D:113:PHE:HD2	1.72	0.55
1:A:23:THR:CG2	1:A:25:GLU:HB3	2.36	0.55
1:A:79:LEU:HD22	1:D:69:GLN:OE1	2.06	0.55
1:B:145:ILE:O	1:B:149:ILE:HG13	2.07	0.55
1:A:82:PRO:O	1:A:86:ILE:HG13	2.07	0.55
1:A:157:LEU:HD22	3:A:428:HOH:O	2.06	0.55
1:B:104:ILE:O	1:B:107:VAL:HB	2.07	0.55
1:C:129:TYR:C	1:C:129:TYR:HD1	2.15	0.55
1:D:8:ASN:O	1:D:8:ASN:ND2	2.40	0.55
1:A:73:LEU:HD22	1:D:73:LEU:HD11	1.88	0.54
1:B:141:PHE:O	1:B:145:ILE:HG23	2.07	0.54
1:A:58:ARG:NH1	1:D:93:LEU:HD13	2.23	0.54
1:A:113:PHE:CE2	1:B:102:LEU:HD22	2.42	0.54
1:A:94:TYR:HA	1:D:62:ASN:ND2	2.14	0.54
1:A:117:GLY:CA	1:B:114:THR:HA	2.37	0.54
1:D:79:LEU:O	1:D:84:GLN:HB2	2.07	0.54
1:C:123:LYS:N	1:C:124:PRO:HD3	2.23	0.54
1:D:27:ILE:HG21	1:D:51:VAL:HG13	1.89	0.54
1:B:94:TYR:CZ	1:B:96:VAL:HB	2.44	0.53
1:D:23:THR:C	1:D:25:GLU:N	2.64	0.53
1:D:35:GLN:OE1	1:D:154:SER:HA	2.09	0.53
1:D:89:ARG:O	1:D:90:ASP:C	2.51	0.53
1:A:130:LEU:HB3	1:A:143:ASP:HB3	1.90	0.53
1:A:108:TYR:OH	2:A:401:SO4:O3	2.27	0.53
1:D:11:GLN:O	1:D:13:ALA:N	2.42	0.53
1:A:160:ARG:HD3	3:A:413:HOH:O	2.07	0.53
1:C:124:PRO:O	1:C:127:LEU:HB2	2.09	0.53
1:D:42:LEU:HD11	1:D:46:GLU:OE2	2.09	0.53
1:A:79:LEU:O	1:A:84:GLN:HB2	2.09	0.53
1:C:72:LYS:NZ	3:C:426:HOH:O	2.40	0.53
1:D:9:LEU:HD12	1:D:9:LEU:N	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:ILE:HG22	1:A:68:ILE:O	2.09	0.52
1:C:110:SER:O	1:C:113:PHE:HD2	1.91	0.52
1:B:58:ARG:HD3	1:C:93:LEU:HB3	1.91	0.52
1:C:10:GLU:OE2	1:C:56:SER:HB3	2.09	0.52
1:A:62:ASN:ND2	1:D:94:TYR:HA	2.24	0.52
1:A:107:VAL:HG12	1:A:108:TYR:CD1	2.45	0.52
1:D:102:LEU:O	1:D:106:ASN:HB2	2.09	0.52
1:A:69:GLN:HE21	1:A:69:GLN:CA	2.17	0.52
1:A:121:LYS:HE2	1:B:115:ASN:ND2	2.25	0.52
1:A:20:ARG:HG3	1:A:68:ILE:HD13	1.92	0.52
1:C:117:GLY:O	1:C:120:ASP:HB3	2.10	0.52
1:B:81:GLU:OE2	1:B:81:GLU:HA	2.10	0.51
1:B:86:ILE:HA	1:B:91:GLU:HB2	1.92	0.51
1:C:96:VAL:O	1:C:99:ILE:HB	2.11	0.51
1:C:114:THR:HG22	1:D:121:LYS:HB2	1.90	0.51
1:D:69:GLN:O	1:D:73:LEU:HB2	2.10	0.51
1:A:157:LEU:HD12	1:A:157:LEU:O	2.10	0.51
1:A:35:GLN:HG2	1:A:154:SER:OG	2.11	0.51
1:B:9:LEU:HD21	1:C:86:ILE:HD11	1.92	0.51
1:A:94:TYR:CA	1:D:62:ASN:HD21	2.16	0.51
1:B:64:ILE:HG12	1:B:145:ILE:HD13	1.91	0.51
1:D:139:ASN:HA	1:D:142:LEU:CD2	2.38	0.51
1:A:83:LEU:HD12	1:D:69:GLN:HB2	1.92	0.51
1:B:76:ASP:O	1:B:78:ARG:HG3	2.11	0.51
1:D:132:ASP:C	1:D:134:SER:N	2.68	0.51
1:A:83:LEU:HG	1:D:65:LEU:HB3	1.92	0.51
1:C:5:ILE:HG23	1:C:6:MSE:N	2.24	0.51
1:A:78:ARG:HD2	1:D:78:ARG:CD	2.41	0.50
1:C:53:ARG:HG3	1:C:156:ARG:CZ	2.41	0.50
1:D:9:LEU:O	1:D:11:GLN:N	2.44	0.50
1:C:110:SER:HA	1:C:113:PHE:CD2	2.47	0.50
1:D:98:GLU:HG2	3:D:409:HOH:O	2.09	0.50
1:A:121:LYS:HE2	1:B:115:ASN:HD21	1.76	0.50
1:A:82:PRO:HD2	1:D:12:THR:HG21	1.93	0.50
1:A:109:GLY:HA3	3:C:432:HOH:O	2.12	0.50
1:A:34:LEU:HD11	1:A:119:ILE:HD13	1.94	0.50
1:C:116:TYR:CE1	1:D:114:THR:HG22	2.47	0.50
1:D:105:VAL:HG23	1:D:148:ALA:HB1	1.94	0.50
1:D:122:GLN:O	1:D:123:LYS:HB2	2.12	0.49
1:B:49:GLU:O	1:B:53:ARG:HG2	2.12	0.49
1:A:142:LEU:HA	1:A:145:ILE:HG12	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:LEU:HD12	1:A:157:LEU:C	2.38	0.49
1:B:111:ILE:HD12	1:B:111:ILE:O	2.12	0.49
1:A:142:LEU:HD12	1:A:145:ILE:CG1	2.43	0.49
1:D:63:ALA:O	1:D:66:THR:N	2.45	0.49
1:B:124:PRO:O	1:B:127:LEU:HB2	2.13	0.49
1:A:158:ALA:O	1:C:4:PHE:HE2	1.95	0.49
1:B:108:TYR:CZ	1:B:156:ARG:HD2	2.47	0.48
1:A:31:VAL:HG11	3:A:428:HOH:O	2.13	0.48
1:B:60:VAL:O	1:B:64:ILE:HG13	2.13	0.48
1:B:131:ASN:O	1:B:132:ASP:C	2.56	0.48
1:D:51:VAL:O	1:D:55:ILE:HG13	2.12	0.48
1:A:126:ILE:O	1:A:129:TYR:N	2.45	0.48
1:D:24:VAL:HG11	1:D:52:ASN:OD1	2.12	0.48
1:A:69:GLN:HB2	1:D:83:LEU:HD12	1.96	0.48
1:D:23:THR:CG2	1:D:26:LYS:HD3	2.43	0.48
1:C:145:ILE:O	1:C:149:ILE:HG13	2.13	0.48
1:D:23:THR:HG23	1:D:26:LYS:H	1.78	0.48
1:C:2:SER:C	1:C:4:PHE:H	2.22	0.48
1:C:113:PHE:CE2	1:D:102:LEU:HD11	2.49	0.48
1:A:15:ARG:HG2	1:A:19:GLU:OE2	2.13	0.48
1:D:134:SER:C	1:D:136:GLY:N	2.72	0.48
1:D:133:LYS:O	1:D:134:SER:HB2	2.13	0.47
1:A:26:LYS:O	1:A:29:GLU:HB2	2.13	0.47
1:A:73:LEU:HB3	1:A:79:LEU:HG	1.96	0.47
1:D:142:LEU:HA	1:D:145:ILE:CD1	2.43	0.47
1:D:39:HIS:N	1:D:40:PRO:HD3	2.30	0.47
1:B:11:GLN:O	1:B:15:ARG:HG3	2.15	0.47
1:D:88:ARG:HG2	1:D:88:ARG:HH11	1.80	0.47
1:B:145:ILE:HD12	1:B:149:ILE:HD11	1.97	0.47
1:D:143:ASP:OD1	1:D:143:ASP:O	2.32	0.47
1:C:34:LEU:CD2	1:C:119:ILE:HG23	2.43	0.47
1:C:156:ARG:NH1	1:C:160:ARG:HD3	2.30	0.47
1:C:156:ARG:HH22	1:C:160:ARG:HH11	1.62	0.47
1:B:101:ALA:HB1	1:B:148:ALA:HB2	1.97	0.46
1:A:118:TYR:HB2	1:B:121:LYS:HG3	1.97	0.46
1:A:23:THR:HG21	1:A:25:GLU:HB3	1.97	0.46
1:C:5:ILE:O	1:C:6:MSE:C	2.57	0.46
1:B:83:LEU:HD22	1:B:87:ILE:HG13	1.98	0.46
1:D:55:ILE:HG12	1:D:149:ILE:HD13	1.96	0.46
1:D:116:TYR:HD2	1:D:148:ALA:HA	1.81	0.46
1:D:70:LEU:HD12	1:D:141:PHE:CZ	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:GLY:HA3	1:B:114:THR:HA	1.96	0.46
1:D:49[A]:GLU:O	1:D:53:ARG:HG2	2.15	0.46
1:D:110:SER:HA	1:D:113:PHE:HE2	1.79	0.46
1:B:129:TYR:CE1	1:B:135:THR:HG21	2.50	0.46
1:D:23:THR:O	1:D:25:GLU:N	2.48	0.46
1:D:25:GLU:HB2	3:D:416:HOH:O	2.15	0.46
1:A:23:THR:HG22	1:A:25:GLU:HB3	1.98	0.46
1:A:40:PRO:CG	1:A:41:ASP:H	2.28	0.46
1:A:34:LEU:HD21	1:A:119:ILE:HG23	1.98	0.45
1:D:30:LEU:HD11	1:D:127:LEU:HG	1.98	0.45
1:D:46:GLU:O	1:D:47:CYS:C	2.59	0.45
1:B:23:THR:CG2	1:B:25:GLU:HB2	2.46	0.45
1:B:77:GLY:HA2	1:B:84:GLN:CD	2.41	0.45
1:C:139:ASN:O	1:C:143:ASP:HB2	2.16	0.45
1:D:22:VAL:CG2	1:D:130:LEU:HD11	2.45	0.45
1:D:142:LEU:HD23	1:D:142:LEU:H	1.80	0.45
1:B:126:ILE:O	1:B:129:TYR:N	2.50	0.45
1:C:81:GLU:OE2	1:C:81:GLU:HA	2.17	0.45
1:A:117:GLY:HA2	1:B:114:THR:HA	1.98	0.45
1:C:27:ILE:HG12	1:C:146:VAL:HG13	1.98	0.45
1:A:101:ALA:HB1	1:A:148:ALA:HB2	1.99	0.45
1:B:89[B]:ARG:O	1:B:90:ASP:HB3	2.16	0.45
1:B:126:ILE:HD12	1:B:126:ILE:N	2.32	0.45
1:D:139:ASN:OD1	1:D:142:LEU:HD11	2.16	0.45
1:C:126:ILE:N	1:C:126:ILE:HD12	2.32	0.45
1:A:125:GLY:HA3	3:A:427:HOH:O	2.17	0.45
1:C:144:ASP:OD2	1:C:144:ASP:N	2.50	0.45
1:A:43:THR:HG23	1:A:46:GLU:HB2	1.99	0.45
1:A:57:LYS:HE2	1:A:108:TYR:OH	2.17	0.45
1:C:13:ALA:HB2	1:C:61:GLN:HG2	1.98	0.45
1:A:17:LEU:HD23	1:A:68:ILE:HD11	1.98	0.44
1:A:70:LEU:HD21	1:D:70:LEU:CD2	2.31	0.44
1:A:115:ASN:O	1:A:119:ILE:HG12	2.16	0.44
1:B:89[B]:ARG:NH1	1:C:3:GLU:HB2	2.32	0.44
1:B:126:ILE:O	1:B:127:LEU:C	2.61	0.44
1:A:160:ARG:O	1:A:161:ALA:C	2.60	0.44
1:A:38:TYR:N	1:A:38:TYR:CD1	2.84	0.44
1:B:93:LEU:HD21	1:C:4:PHE:HE2	1.80	0.44
1:C:32:TYR:O	1:C:36:SER:HB2	2.17	0.44
1:B:158:ALA:O	1:B:159:HIS:C	2.58	0.44
1:D:60:VAL:HG23	1:D:107:VAL:HG21	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:30:LEU:HD11	1:B:126:ILE:HD11	1.99	0.44
1:D:101:ALA:HB2	1:D:144:ASP:HB3	1.99	0.44
1:A:26:LYS:HB3	1:A:126:ILE:HG12	1.99	0.44
1:A:126:ILE:O	1:A:127:LEU:C	2.61	0.44
1:D:88:ARG:HG2	1:D:88:ARG:NH1	2.32	0.44
1:D:108:TYR:CZ	1:D:156:ARG:HD2	2.52	0.44
1:A:26:LYS:HG2	1:A:29:GLU:OE2	2.18	0.43
1:B:98:GLU:HG3	1:B:144:ASP:OD1	2.18	0.43
1:C:11:GLN:HG2	1:C:15:ARG:NH2	2.33	0.43
1:C:63:ALA:HA	1:C:100:LEU:HD22	2.00	0.43
1:A:28:ALA:HB3	1:A:44:MSE:HE3	1.99	0.43
1:A:53:ARG:HD2	1:A:53:ARG:HA	1.76	0.43
1:D:10:GLU:HA	1:D:61:GLN:NE2	2.34	0.43
1:A:29:GLU:N	1:A:44:MSE:HE1	2.34	0.43
1:B:30:LEU:HD11	1:B:126:ILE:CD1	2.48	0.43
1:C:4:PHE:O	1:C:5:ILE:C	2.61	0.43
1:A:130:LEU:O	1:A:139:ASN:HB2	2.18	0.43
1:C:132:ASP:HB3	1:C:135:THR:HG23	2.01	0.43
1:B:63:ALA:HB1	1:B:100:LEU:HD12	2.00	0.43
1:C:135:THR:OG1	1:C:137:LYS:HB2	2.19	0.43
1:C:158:ALA:O	1:C:159:HIS:C	2.60	0.43
1:A:100:LEU:HD21	1:D:96:VAL:HG13	2.00	0.43
1:A:159:HIS:HE1	3:B:411:HOH:O	2.01	0.43
1:B:48:ILE:HG22	1:B:52:ASN:ND2	2.33	0.43
1:D:53:ARG:HG3	1:D:156:ARG:CZ	2.48	0.43
1:B:6:MSE:HG3	3:C:428:HOH:O	2.19	0.43
1:B:129:TYR:HE1	1:B:135:THR:HG21	1.84	0.43
1:A:36:SER:O	1:A:37:LYS:C	2.61	0.42
1:D:53:ARG:O	1:D:56:SER:OG	2.31	0.42
1:A:27:ILE:O	1:A:30:LEU:HB2	2.19	0.42
1:B:126:ILE:HD12	1:B:127:LEU:H	1.84	0.42
1:C:11:GLN:HE21	1:C:11:GLN:HB2	1.67	0.42
1:D:63:ALA:O	1:D:64:ILE:C	2.61	0.42
1:A:83:LEU:HD11	1:D:66:THR:CA	2.48	0.42
1:C:88:ARG:HA	1:C:88:ARG:HD2	1.91	0.42
1:A:16:TRP:O	1:A:20:ARG:HG2	2.19	0.42
1:C:107:VAL:HG12	1:C:108:TYR:CD1	2.54	0.42
1:D:9:LEU:C	1:D:11:GLN:N	2.77	0.42
1:D:54:VAL:CG1	1:D:149:ILE:HG12	2.49	0.42
1:D:79:LEU:HD13	1:D:87:ILE:HD12	2.00	0.42
1:B:23:THR:HB	1:B:26:LYS:HG2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:132:ASP:O	1:C:133:LYS:CB	2.59	0.42
1:A:78:ARG:HG3	1:D:78:ARG:HH11	1.85	0.42
1:B:14:ARG:O	1:B:18:GLU:HG3	2.20	0.42
1:C:54:VAL:HG22	1:C:108:TYR:CE1	2.55	0.42
1:B:118:TYR:CE2	1:B:122:GLN:HG3	2.55	0.42
1:C:32:TYR:CZ	1:C:36:SER:OG	2.69	0.42
1:B:63:ALA:HA	1:B:100:LEU:CD1	2.49	0.42
1:C:43:THR:HG1	1:C:46:GLU:HG2	1.80	0.42
1:D:142:LEU:HA	1:D:145:ILE:HD11	2.02	0.42
1:A:142:LEU:CD1	1:A:145:ILE:HD11	2.49	0.41
1:B:33:TYR:C	1:B:33:TYR:HD2	2.26	0.41
1:C:5:ILE:HG23	1:C:6:MSE:H	1.85	0.41
1:A:130:LEU:N	1:A:130:LEU:HD12	2.35	0.41
1:B:119:ILE:C	1:B:121:LYS:H	2.28	0.41
1:A:-2:SER:O	1:A:2:SER:HB2	2.21	0.41
1:B:85:SER:O	1:B:89[A]:ARG:HB2	2.20	0.41
1:C:72:LYS:HE2	3:C:435:HOH:O	2.21	0.41
1:D:34:LEU:HD11	1:D:119:ILE:CG2	2.50	0.41
1:C:20:ARG:NH2	1:C:75:GLU:OE2	2.53	0.41
1:D:35:GLN:NE2	1:D:157:LEU:HD22	2.34	0.41
1:B:23:THR:HG22	1:B:25:GLU:HB2	2.03	0.41
1:B:48:ILE:HG22	1:B:52:ASN:HD21	1.86	0.41
1:C:15:ARG:NH1	3:C:425:HOH:O	2.53	0.41
1:D:24:VAL:HG13	1:D:51:VAL:HG12	2.02	0.41
1:A:4:PHE:O	1:A:8:ASN:ND2	2.42	0.41
1:C:28:ALA:HA	1:C:51:VAL:HG21	2.03	0.41
1:C:57:LYS:HE2	1:C:108:TYR:OH	2.21	0.41
1:C:116:TYR:O	1:C:120:ASP:HB2	2.20	0.41
1:D:60:VAL:O	1:D:64:ILE:HG13	2.20	0.41
1:A:89:ARG:O	1:A:90:ASP:HB3	2.21	0.41
1:A:78:ARG:CZ	1:D:78:ARG:HD2	2.50	0.40
1:B:54:VAL:HG11	1:B:149:ILE:HG23	2.02	0.40
1:B:84:GLN:HE21	1:B:88[B]:ARG:HD2	1.85	0.40
1:A:2:SER:OG	1:D:91:GLU:OE2	2.28	0.40
1:A:40:PRO:HG2	1:A:41:ASP:N	2.30	0.40
1:A:142:LEU:HD12	1:A:145:ILE:CD1	2.50	0.40
1:D:116:TYR:O	1:D:120:ASP:HB2	2.20	0.40
1:A:77:GLY:HA2	1:A:84:GLN:NE2	2.37	0.40
1:B:89[B]:ARG:HB3	1:B:89[B]:ARG:NH1	2.34	0.40
1:D:156:ARG:O	1:D:159:HIS:N	2.55	0.40
1:B:23:THR:CG2	1:B:24:VAL:N	2.84	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:25:GLU:O	1:B:28:ALA:HB3	2.20	0.40
1:C:60:VAL:O	1:C:61:GLN:C	2.62	0.40
1:A:22:VAL:CG2	1:A:142:LEU:HD21	2.51	0.40
1:D:9:LEU:N	1:D:9:LEU:CD1	2.85	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	162/168 (96%)	149 (92%)	13 (8%)	0	100	100
1	B	164/168 (98%)	149 (91%)	13 (8%)	2 (1%)	10	34
1	C	159/168 (95%)	140 (88%)	16 (10%)	3 (2%)	6	22
1	D	153/168 (91%)	123 (80%)	24 (16%)	6 (4%)	2	8
All	All	638/672 (95%)	561 (88%)	66 (10%)	11 (2%)	7	25

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	5	ILE
1	C	40	PRO
1	B	132	ASP
1	D	12	THR
1	D	132	ASP
1	D	11	GLN
1	C	161	ALA
1	D	10	GLU
1	D	64	ILE
1	D	24	VAL
1	B	125	GLY

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	140/139 (101%)	126 (90%)	14 (10%)	7	24
1	B	142/139 (102%)	129 (91%)	13 (9%)	8	27
1	C	137/139 (99%)	126 (92%)	11 (8%)	11	34
1	D	132/139 (95%)	122 (92%)	10 (8%)	12	36
All	All	551/556 (99%)	503 (91%)	48 (9%)	9	30

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	VAL
1	A	25	GLU
1	A	43	THR
1	A	49	GLU
1	A	59	GLU
1	A	69	GLN
1	A	91	GLU
1	A	114	THR
1	A	121	LYS
1	A	122	GLN
1	A	126	ILE
1	A	131	ASN
1	A	132	ASP
1	A	157	LEU
1	B	-2	SER
1	B	33	TYR
1	B	83	LEU
1	B	98	GLU
1	B	100	LEU
1	B	102	LEU
1	B	107	VAL
1	B	111	ILE
1	B	114	THR
1	B	115	ASN

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Mol	Chain	Res	Type
1	B	126	ILE
1	B	133	LYS
1	B	145	ILE
1	C	4	PHE
1	C	8	ASN
1	C	20	ARG
1	C	49	GLU
1	C	58	ARG
1	C	83	LEU
1	C	88	ARG
1	C	111	ILE
1	C	119	ILE
1	C	126	ILE
1	C	129	TYR
1	D	20	ARG
1	D	25	GLU
1	D	34	LEU
1	D	69	GLN
1	D	78	ARG
1	D	83	LEU
1	D	98	GLU
1	D	132	ASP
1	D	142	LEU
1	D	149	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	HIS
1	A	122	GLN
1	A	128	GLN
1	A	159	HIS
1	B	8	ASN
1	B	35	GLN
1	B	115	ASN
1	B	122	GLN
1	B	131	ASN
1	B	139	ASN
1	C	8	ASN
1	C	122	GLN
1	D	8	ASN
1	D	39	HIS

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Mol	Chain	Res	Type
1	D	62	ASN
1	D	106	ASN
1	D	122	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](#)

7 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	A	406	-	4,4,4	0.55	0	6,6,6	0.22	0
2	SO4	C	403	-	4,4,4	0.41	0	6,6,6	0.14	0
2	SO4	C	407	-	4,4,4	0.74	0	6,6,6	0.15	0
2	SO4	B	402	-	4,4,4	0.43	0	6,6,6	0.13	0
2	SO4	D	404	-	4,4,4	0.78	0	6,6,6	0.12	0
2	SO4	A	401	-	4,4,4	0.47	0	6,6,6	0.18	0
2	SO4	A	405	-	4,4,4	0.74	0	6,6,6	0.11	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	404	SO4	1	0
2	A	401	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	161/168 (95%)	0.30	2 (1%) 76 68	40, 67, 98, 123	0
1	B	161/168 (95%)	-0.17	3 (1%) 66 57	20, 44, 78, 106	2 (1%)
1	C	159/168 (94%)	-0.05	8 (5%) 34 26	27, 46, 87, 127	0
1	D	153/168 (91%)	0.84	12 (7%) 19 14	40, 90, 117, 132	1 (0%)
All	All	634/672 (94%)	0.22	25 (3%) 43 34	20, 63, 105, 132	3 (0%)

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	161	ALA	5.8
1	C	4	PHE	4.8
1	A	161	ALA	3.0
1	C	162	ALA	2.9
1	C	5	ILE	2.8
1	B	109	GLY	2.7
1	D	142	LEU	2.7
1	B	159	HIS	2.6
1	D	143	ASP	2.5
1	C	129	TYR	2.5
1	A	86	ILE	2.4
1	C	41	ASP	2.3
1	D	108	TYR	2.3
1	D	129	TYR	2.3
1	B	161	ALA	2.2
1	D	74	ALA	2.2
1	D	136	GLY	2.1
1	D	132	ASP	2.1
1	D	9	LEU	2.1
1	C	161	ALA	2.1
1	D	113	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	37	LYS	2.0
1	D	150	ALA	2.0
1	D	112	GLY	2.0
1	C	111	ILE	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($\text{\AA}^2$)	Q<0.9
2	SO4	D	404	5/5	0.80	0.14	77,77,78,82	0
2	SO4	C	407	5/5	0.84	0.18	77,78,79,82	0
2	SO4	A	406	5/5	0.87	0.17	69,69,72,73	0
2	SO4	A	405	5/5	0.90	0.15	74,75,78,81	0
2	SO4	A	401	5/5	0.90	0.08	65,66,68,68	0
2	SO4	C	403	5/5	0.93	0.07	63,64,66,67	0
2	SO4	B	402	5/5	0.94	0.07	61,62,64,65	0

6.5 Other polymers [i](#)

There are no such residues in this entry.