



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:23 AM UTC

PDB ID : 4RJY / pdb_00004rjy
Title : Crystal structure of E. coli L-Threonine Aldolase in complex with a non-covalently uncleaved bound L-serine substrate
Authors : Safo, M.K.; Chowdhury, N.; Gandhi, A.K.
Deposited on : 2014-10-11
Resolution : 2.10 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

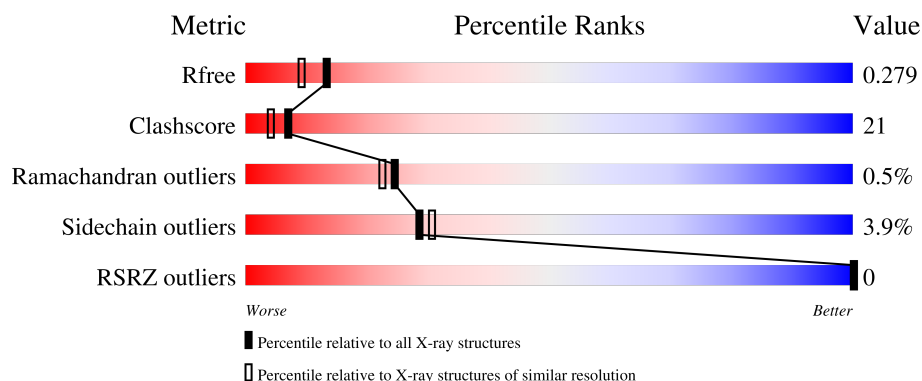
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	333	69% 28% .
1	B	333	62% 33% 5%
1	C	333	53% 41% 5%
1	D	333	65% 33% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11165 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 Low specificity L-threonine aldolase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	332	Total	C	N	O	P	S	0	0	0
			2563	1606	459	481	1	16			
1	B	332	Total	C	N	O	P	S	0	0	0
			2563	1606	459	481	1	16			
1	C	332	Total	C	N	O	P	S	0	0	0
			2563	1606	459	481	1	16			
1	D	332	Total	C	N	O	P	S	0	0	0
			2563	1606	459	481	1	16			

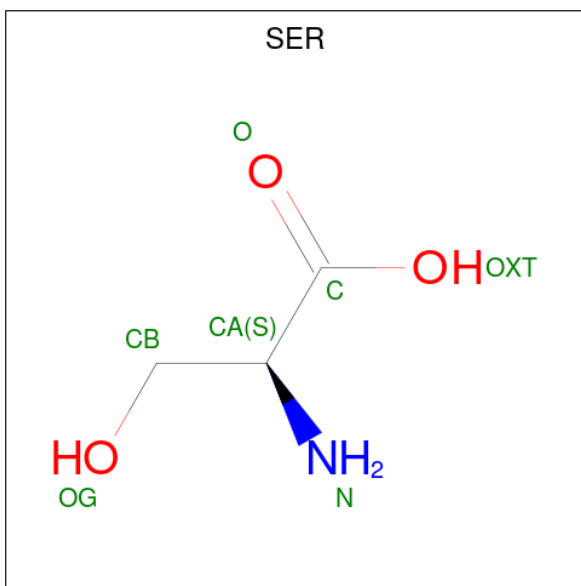
There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	158	LYS	GLU	conflict	UNP L4EJM2
A	256	THR	ALA	conflict	UNP L4EJM2
B	158	LYS	GLU	conflict	UNP L4EJM2
B	256	THR	ALA	conflict	UNP L4EJM2
C	158	LYS	GLU	conflict	UNP L4EJM2
C	256	THR	ALA	conflict	UNP L4EJM2
D	158	LYS	GLU	conflict	UNP L4EJM2
D	256	THR	ALA	conflict	UNP L4EJM2

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Na	0	0
			1	1		
2	B	2	Total	Na	0	0
			2	2		
2	C	1	Total	Na	0	0
			1	1		
2	D	2	Total	Na	0	0
			2	2		

- Molecule 3 is SERINE (CCD ID: SER) (formula: $C_3H_7NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	1
			14	6	2	6		
3	B	1	Total	C	N	O	0	1
			14	6	2	6		
3	C	1	Total	C	N	O	0	1
			14	6	2	6		
3	D	1	Total	C	N	O	0	1
			14	6	2	6		

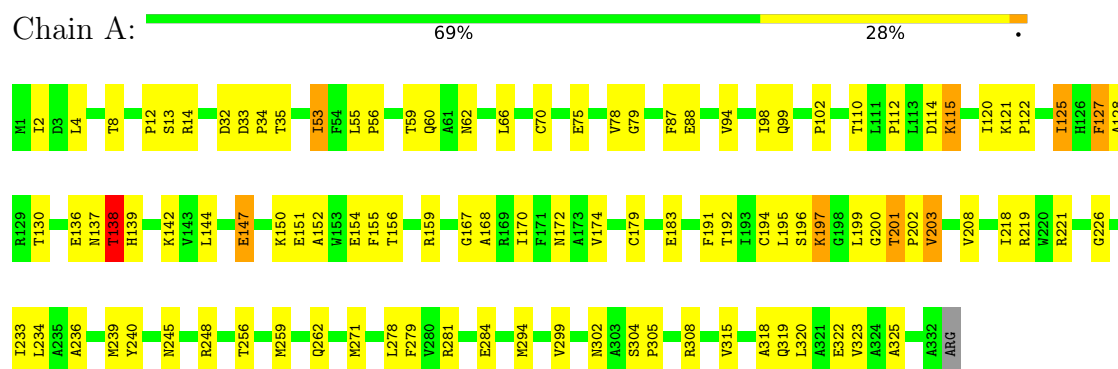
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	230	Total	O	0	0
			230	230		
4	B	194	Total	O	0	0
			194	194		
4	C	207	Total	O	0	0
			207	207		
4	D	220	Total	O	0	0
			220	220		

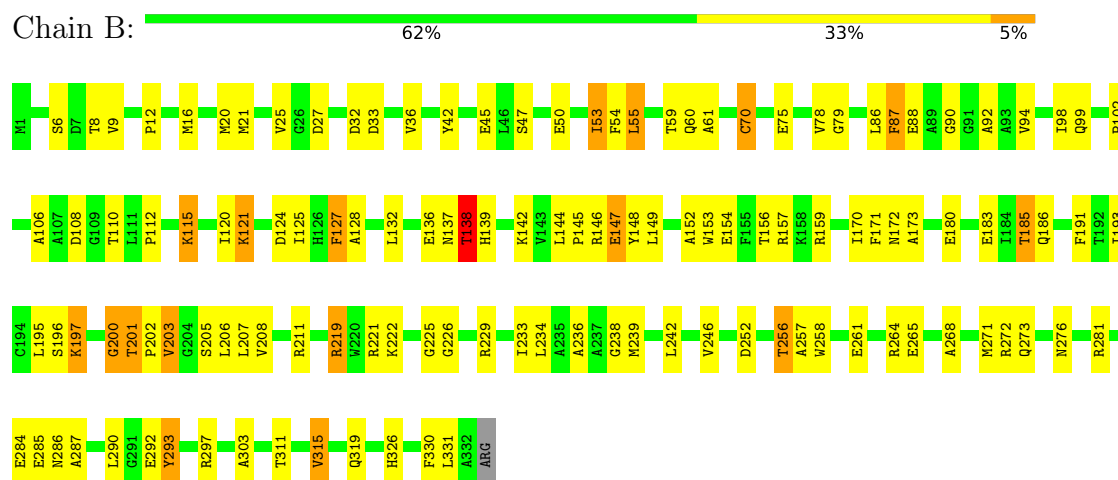
3 Residue-property plots

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.

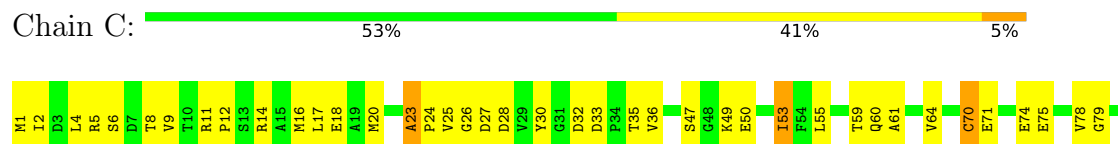
- Molecule 1: Low specificity L-threonine aldolase

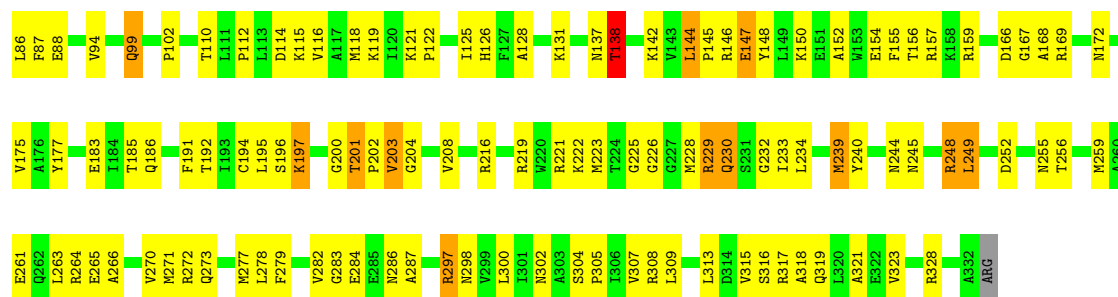


- Molecule 1: Low specificity L-threonine aldolase



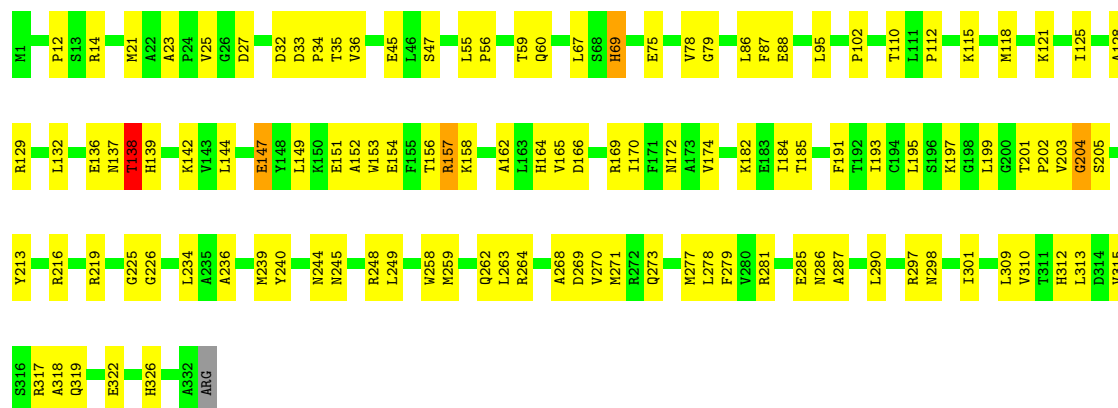
- Molecule 1: Low specificity L-threonine aldolase





- Molecule 1: Low specificity L-threonine aldolase

Chain D: 65% 33% .



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	77.16Å 104.88Å 84.91Å 90.00° 92.49° 90.00°	Depositor
Resolution (Å)	30.00 – 2.10 30.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.6 (30.00-2.10) 97.6 (30.00-2.10)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.07 (at 2.10Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.212 , 0.279 0.214 , 0.279	Depositor DCC
R_{free} test set	3884 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	21.6	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.056 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11165	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.84% 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: NA, LLP

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	0.55	0/2582	1.04	10/3496 (0.3%)
1	B	0.54	0/2582	1.03	15/3496 (0.4%)
1	C	0.54	0/2582	1.06	12/3496 (0.3%)
1	D	0.54	0/2582	1.05	11/3496 (0.3%)
All	All	0.55	0/10328	1.05	48/13984 (0.3%)

There are no bond length outliers.

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	138	THR	N-CA-C	-9.58	90.40	110.80
1	A	138	THR	N-CA-C	-9.16	91.28	110.80
1	C	138	THR	N-CA-C	-8.87	91.90	110.80
1	D	87	PHE	N-CA-C	8.01	122.77	112.92
1	B	138	THR	N-CA-C	-7.65	94.50	110.80
1	C	128	ALA	N-CA-C	-7.58	100.52	110.53
1	B	172	ASN	N-CA-C	-7.38	103.23	111.71
1	B	121	LYS	N-CA-C	7.37	119.06	109.93
1	D	128	ALA	N-CA-C	-7.02	101.07	110.55
1	C	200	GLY	N-CA-C	7.01	124.66	115.47
1	B	55	LEU	CA-C-N	6.95	126.76	119.05
1	B	55	LEU	C-N-CA	6.95	126.76	119.05
1	B	70	CYS	N-CA-C	6.61	119.70	109.07
1	B	200	GLY	N-CA-C	6.48	125.24	114.48
1	C	204	GLY	N-CA-C	6.44	125.47	111.04
1	D	201	THR	N-CA-C	-6.32	102.73	110.31
1	C	23	ALA	N-CA-C	6.14	117.43	109.64
1	D	25	VAL	N-CA-C	6.06	116.22	108.82
1	C	172	ASN	N-CA-C	-6.03	104.83	111.82
1	A	87	PHE	N-CA-C	6.00	119.64	112.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	245	ASN	N-CA-C	5.99	120.28	112.92
1	B	246	VAL	N-CA-C	5.98	116.18	110.74
1	B	128	ALA	N-CA-C	-5.94	102.69	110.53
1	C	249	LEU	N-CA-C	-5.85	105.03	111.82
1	B	257	ALA	N-CA-C	-5.77	105.07	111.36
1	A	172	ASN	N-CA-C	-5.74	105.02	111.28
1	D	204	GLY	N-CA-C	5.74	126.78	113.18
1	D	172	ASN	N-CA-C	-5.66	105.26	111.82
1	C	70	CYS	N-CA-C	5.64	118.15	109.07
1	C	87	PHE	N-CA-C	5.64	119.25	112.93
1	A	200	GLY	N-CA-C	5.57	123.03	115.30
1	B	106	ALA	N-CA-C	-5.42	103.54	110.53
1	B	127	PHE	N-CA-C	5.38	118.48	110.14
1	C	94	VAL	N-CA-C	5.38	115.58	110.42
1	A	33	ASP	CA-C-N	5.31	125.12	119.28
1	A	33	ASP	C-N-CA	5.31	125.12	119.28
1	A	94	VAL	N-CA-C	5.26	115.48	110.53
1	C	30	TYR	N-CA-C	-5.25	105.62	112.23
1	A	127	PHE	N-CA-C	5.24	118.28	109.95
1	D	244	ASN	N-CA-C	5.23	120.66	113.97
1	C	282	VAL	N-CA-C	5.21	115.87	110.82
1	B	25	VAL	N-CA-C	5.20	115.66	108.23
1	A	13	SER	N-CA-C	-5.20	103.53	110.55
1	D	166	ASP	N-CA-C	-5.19	100.63	108.67
1	B	132	LEU	N-CA-C	5.18	117.69	109.50
1	D	157	ARG	N-CA-C	-5.15	105.74	111.36
1	B	87	PHE	N-CA-C	5.11	118.66	112.93
1	A	125	ILE	N-CA-C	5.08	119.21	111.89

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	2563	0	2552	92	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2563	0	2554	111	0
1	C	2563	0	2553	134	0
1	D	2563	0	2553	111	0
2	A	1	0	0	0	0
2	B	2	0	0	0	0
2	C	1	0	0	0	0
2	D	2	0	0	0	0
3	A	14	0	8	1	0
3	B	14	0	8	1	0
3	C	14	0	8	0	0
3	D	14	0	8	0	0
4	A	230	0	0	14	0
4	B	194	0	0	14	0
4	C	207	0	0	13	0
4	D	220	0	0	15	0
All	All	11165	0	10244	431	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 21.

All (431) 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:236:ALA:HA	1:B:239:MET:HE2	1.39	1.02
1:B:60:GLN:HE22	1:B:226:GLY:HA3	1.25	1.02
1:D:59:THR:HG23	1:D:88:GLU:HG2	1.37	1.02
1:B:12:PRO:HG3	1:B:202:PRO:HG3	1.43	1.01
1:A:60:GLN:HE22	1:A:226:GLY:HA3	1.23	1.00
1:D:236:ALA:HA	1:D:239:MET:HE2	1.43	0.98
1:A:219:ARG:HG2	1:C:125:ILE:HG22	1.44	0.97
1:C:167:GLY:H	1:C:192:THR:HG22	1.30	0.96
1:D:278:LEU:HD11	1:D:309:LEU:HD12	1.49	0.95
1:B:47:SER:HB2	1:B:185:THR:HG21	1.49	0.93
1:A:12:PRO:HG3	1:A:202:PRO:HG3	1.53	0.89
1:A:137:ASN:HD22	1:A:138:THR:N	1.72	0.88
1:B:153:TRP:O	1:B:157:ARG:HG2	1.75	0.86
1:C:60:GLN:HE22	1:C:226:GLY:HA3	1.39	0.86
1:B:55:LEU:HD13	1:B:61:ALA:HA	1.58	0.84
1:A:59:THR:HG23	1:A:88:GLU:HG2	1.60	0.82
1:D:12:PRO:HG3	1:D:202:PRO:HG3	1.58	0.82
1:C:12:PRO:HG3	1:C:202:PRO:HG3	1.60	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:196:SER:HA	1:C:201:THR:HG22	1.65	0.79
1:C:138:THR:HG22	4:C:706:HOH:O	1.83	0.78
1:A:271:MET:HE3	1:A:281:ARG:NH2	1.99	0.77
1:B:59:THR:HG23	1:B:88:GLU:HG2	1.64	0.77
1:C:35:THR:HG23	1:C:239:MET:HE1	1.68	0.76
1:B:264:ARG:HH21	1:B:273:GLN:HE21	1.34	0.75
1:A:138:THR:HA	1:A:142:LYS:O	1.87	0.75
1:C:167:GLY:N	1:C:192:THR:HG22	2.02	0.75
1:A:75:GLU:HG3	1:A:130:THR:HA	1.68	0.75
1:D:60:GLN:HE22	1:D:226:GLY:HA3	1.52	0.75
1:B:256:THR:HG21	1:B:276:ASN:HA	1.68	0.74
1:B:147:GLU:CD	1:B:147:GLU:H	1.96	0.74
1:C:155:PHE:O	1:C:159:ARG:HG3	1.88	0.73
1:D:115:LYS:HA	1:D:118:MET:HE3	1.70	0.73
1:B:219:ARG:NH1	1:D:125:ILE:HA	2.04	0.73
1:B:196:SER:HA	1:B:201:THR:HG22	1.68	0.73
1:D:182:LYS:HA	1:D:185:THR:HG22	1.71	0.73
1:C:78:VAL:O	1:C:102:PRO:HA	1.89	0.72
1:C:47:SER:HB2	1:C:185:THR:HG21	1.71	0.72
1:B:196:SER:CA	1:B:201:THR:HG22	2.19	0.72
1:C:4:LEU:HD21	1:C:315:VAL:HG21	1.71	0.72
1:C:26:GLY:HA3	1:C:32:ASP:O	1.90	0.72
1:C:196:SER:CA	1:C:201:THR:HG22	2.20	0.72
1:A:60:GLN:NE2	1:A:226:GLY:HA3	2.03	0.71
1:B:285:GLU:HG3	1:B:286:ASN:OD1	1.91	0.71
1:D:153:TRP:O	1:D:157:ARG:HG2	1.91	0.71
1:D:55:LEU:HD23	1:D:60:GLN:HE21	1.56	0.71
1:D:147:GLU:CD	1:D:147:GLU:H	1.99	0.71
1:A:203:VAL:HG21	1:A:234:LEU:HD13	1.72	0.70
1:B:315:VAL:HA	1:B:319:GLN:OE1	1.90	0.69
1:A:203:VAL:CG2	1:A:234:LEU:HD13	2.21	0.69
1:D:138:THR:HG21	1:D:169:ARG:HB2	1.75	0.69
1:B:60:GLN:NE2	1:B:226:GLY:HA3	2.05	0.68
1:A:125:ILE:HG22	1:C:219:ARG:HG2	1.75	0.68
1:A:147:GLU:O	1:A:151:GLU:HG3	1.92	0.68
1:C:195:LEU:HB2	1:C:201:THR:HG21	1.75	0.68
1:D:312:HIS:CE1	1:D:315:VAL:HG23	2.28	0.68
1:D:170:ILE:O	1:D:174:VAL:HG23	1.94	0.67
1:C:216:ARG:HD2	4:C:627:HOH:O	1.94	0.66
1:A:78:VAL:HG22	1:A:79:GLY:N	2.10	0.66
1:C:261:GLU:O	1:C:265:GLU:HG3	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:137:ASN:HD22	1:D:138:THR:N	1.94	0.66
1:D:138:THR:HA	1:D:142:LYS:O	1.96	0.66
1:D:301:ILE:C	1:D:301:ILE:HD12	2.21	0.66
1:C:146:ARG:HG3	1:C:177:TYR:CE1	2.32	0.65
1:C:27:ASP:H	1:C:32:ASP:HB2	1.61	0.65
1:A:147:GLU:H	1:A:147:GLU:CD	2.04	0.65
1:D:78:VAL:O	1:D:102:PRO:HA	1.96	0.65
1:D:264:ARG:HH11	1:D:264:ARG:HG2	1.62	0.65
1:D:47:SER:HB2	1:D:185:THR:HG21	1.79	0.65
1:A:201:THR:HG23	1:A:202:PRO:HD2	1.77	0.64
1:C:297:ARG:HG3	1:C:297:ARG:HH11	1.62	0.64
1:D:112:PRO:HB2	1:D:115:LYS:HB2	1.79	0.64
1:C:196:SER:N	1:C:201:THR:HG22	2.11	0.64
1:B:16:MET:HE1	1:B:200:GLY:O	1.97	0.64
1:B:115:LYS:NZ	1:B:115:LYS:HB3	2.11	0.64
1:B:256:THR:HG22	1:B:311:THR:CG2	2.28	0.63
1:B:138:THR:HA	1:B:142:LYS:O	1.98	0.63
1:C:168:ALA:HA	1:C:194:CYS:HB2	1.79	0.63
1:A:137:ASN:HD22	1:A:138:THR:H	1.46	0.63
1:D:32:ASP:O	1:D:34:PRO:HD3	1.98	0.63
1:C:8:THR:HG22	1:C:197:LLP:HG3	1.80	0.62
1:C:270:VAL:HB	4:C:630:HOH:O	1.98	0.62
1:D:149:LEU:HD22	4:D:597:HOH:O	1.99	0.62
1:B:264:ARG:HH21	1:B:273:GLN:NE2	1.97	0.62
1:C:192:THR:HG21	4:C:597:HOH:O	2.00	0.62
1:C:167:GLY:H	1:C:192:THR:CG2	2.08	0.62
1:D:264:ARG:HE	1:D:270:VAL:HB	1.63	0.62
1:A:112:PRO:HB2	1:A:115:LYS:HB2	1.81	0.62
1:D:154:GLU:O	1:D:158:LYS:HG3	1.98	0.62
1:A:294:MET:HB3	1:A:299:VAL:HB	1.80	0.62
1:B:125:ILE:HG22	1:D:219:ARG:HG3	1.81	0.61
1:A:179:CYS:HB2	1:A:183:GLU:OE1	2.00	0.61
1:B:252:ASP:O	1:B:256:THR:HG23	1.99	0.61
1:C:230:GLN:H	1:C:230:GLN:HE21	1.49	0.61
1:D:59:THR:HG23	1:D:88:GLU:CG	2.23	0.61
1:C:150:LYS:HG2	1:C:154:GLU:OE2	2.01	0.60
1:D:318:ALA:HB2	4:D:545:HOH:O	1.99	0.60
1:A:138:THR:HG22	4:A:641:HOH:O	2.01	0.60
1:C:14:ARG:O	1:C:18:GLU:HG3	2.02	0.60
1:A:136:GLU:OE2	1:A:139:HIS:HD2	1.85	0.59
1:C:297:ARG:O	1:C:298:ASN:HB2	2.01	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:285:GLU:C	1:D:286:ASN:HD22	2.11	0.59
1:B:8:THR:HG22	1:B:197:LLP:HE3	1.83	0.59
1:D:136:GLU:OE2	1:D:139:HIS:HD2	1.85	0.59
1:A:75:GLU:CG	1:A:130:THR:HA	2.32	0.59
1:A:271:MET:HE3	1:A:281:ARG:HH22	1.68	0.59
1:C:17:LEU:HB3	1:D:21:MET:HE2	1.85	0.58
1:B:152:ALA:O	1:B:156:THR:HG23	2.04	0.58
3:B:403[A]:SER:N	4:B:666:HOH:O	2.36	0.58
1:B:50:GLU:HG3	4:B:632:HOH:O	2.03	0.58
1:C:259:MET:HE1	1:C:323:VAL:HB	1.84	0.58
1:C:297:ARG:HG3	1:C:297:ARG:NH1	2.15	0.58
1:A:152:ALA:O	1:A:156:THR:HG23	2.03	0.58
1:B:33:ASP:HB3	1:B:36:VAL:HB	1.86	0.58
1:C:115:LYS:HG3	4:C:654:HOH:O	2.04	0.58
1:B:268:ALA:HA	4:B:564:HOH:O	2.03	0.58
1:C:8:THR:HG22	1:C:197:LLP:HE2	1.86	0.57
1:D:174:VAL:HG22	1:D:184:ILE:HD11	1.85	0.57
1:B:112:PRO:HA	4:B:638:HOH:O	2.05	0.57
1:B:138:THR:HG22	4:B:692:HOH:O	2.04	0.57
1:C:60:GLN:HB2	1:D:225:GLY:HA3	1.86	0.57
1:A:60:GLN:HE22	1:A:226:GLY:CA	2.07	0.57
1:C:114:ASP:O	1:C:118:MET:HG3	2.05	0.57
1:B:238:GLY:O	1:B:242:LEU:HG	2.05	0.56
1:C:195:LEU:C	1:C:201:THR:HG22	2.30	0.56
1:A:128:ALA:HA	4:A:515:HOH:O	2.04	0.56
1:A:127:PHE:O	1:C:219:ARG:NH2	2.38	0.56
1:A:170:ILE:O	1:A:174:VAL:HG23	2.04	0.56
1:C:59:THR:HG23	1:C:88:GLU:HG2	1.85	0.56
1:D:78:VAL:HG22	1:D:79:GLY:N	2.19	0.56
1:D:110:THR:HG22	1:D:144:LEU:HG	1.87	0.56
1:B:196:SER:N	1:B:201:THR:HG22	2.21	0.56
1:C:99:GLN:HE21	1:C:99:GLN:HA	1.70	0.56
1:B:287:ALA:O	1:B:303:ALA:HB1	2.04	0.56
1:A:2:ILE:HD13	1:A:319:GLN:HB3	1.88	0.56
1:A:150:LYS:O	1:A:154:GLU:HG3	2.06	0.56
1:C:230:GLN:H	1:C:230:GLN:NE2	2.02	0.56
1:A:60:GLN:HB2	1:B:225:GLY:HA3	1.88	0.55
1:A:78:VAL:O	1:A:102:PRO:HA	2.05	0.55
1:C:24:PRO:HD2	1:C:35:THR:OG1	2.06	0.55
1:C:137:ASN:HD21	1:C:144:LEU:H	1.54	0.55
1:C:271:MET:O	1:C:272:ARG:HB3	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:16:MET:O	1:C:20:MET:HG3	2.07	0.55
1:C:55:LEU:HD22	1:C:61:ALA:HA	1.88	0.55
1:C:138:THR:HA	1:C:142:LYS:O	2.06	0.55
1:A:150:LYS:HE3	4:A:591:HOH:O	2.07	0.55
1:A:12:PRO:CG	1:A:202:PRO:HG3	2.32	0.55
1:B:6:SER:O	1:B:9:VAL:HG22	2.06	0.55
1:B:92:ALA:HB3	4:B:532:HOH:O	2.07	0.55
1:D:203:VAL:HG12	1:D:204:GLY:N	2.22	0.55
1:A:78:VAL:HG22	1:A:79:GLY:H	1.70	0.54
1:B:138:THR:HG22	1:B:138:THR:O	2.07	0.54
1:B:219:ARG:HG2	4:B:645:HOH:O	2.06	0.54
1:C:229:ARG:HB3	1:C:230:GLN:NE2	2.22	0.54
1:C:6:SER:O	1:C:9:VAL:HG22	2.07	0.54
1:A:318:ALA:HB2	4:A:655:HOH:O	2.07	0.54
1:D:138:THR:CG2	1:D:169:ARG:HB2	2.36	0.54
1:B:271:MET:HE3	1:B:281:ARG:NH2	2.23	0.54
1:A:8:THR:HG22	1:A:197:LLP:HG3	1.90	0.54
1:C:222:LYS:O	4:C:525:HOH:O	2.18	0.54
1:A:218:ILE:HD11	4:A:688:HOH:O	2.08	0.53
1:C:11:ARG:NH1	4:C:513:HOH:O	2.41	0.53
1:C:35:THR:CG2	1:C:239:MET:HE1	2.37	0.53
1:D:277:MET:HG2	1:D:310:VAL:HG22	1.89	0.53
1:B:137:ASN:O	1:B:138:THR:HB	2.08	0.53
1:B:78:VAL:HG22	1:B:79:GLY:N	2.24	0.53
1:C:8:THR:CG2	1:C:197:LLP:HE2	2.39	0.53
1:D:219:ARG:NH1	4:D:578:HOH:O	2.41	0.53
1:A:53:ILE:HD11	1:A:221:ARG:HD3	1.91	0.53
1:B:138:THR:HG21	4:B:504:HOH:O	2.08	0.53
1:C:230:GLN:NE2	1:C:230:GLN:N	2.57	0.53
1:D:301:ILE:HD12	1:D:301:ILE:O	2.08	0.53
1:B:256:THR:HG21	1:B:276:ASN:CA	2.37	0.53
1:C:53:ILE:HD11	1:C:221:ARG:HD3	1.91	0.53
1:A:256:THR:HG23	1:A:278:LEU:HD22	1.90	0.52
1:C:264:ARG:HD3	4:C:630:HOH:O	2.10	0.52
1:B:264:ARG:NH2	1:B:273:GLN:HE21	2.07	0.52
1:D:137:ASN:O	1:D:138:THR:HB	2.09	0.52
1:B:211:ARG:NH1	1:B:211:ARG:HB2	2.25	0.52
1:A:167:GLY:H	1:A:192:THR:CG2	2.23	0.52
1:C:313:LEU:C	1:C:313:LEU:HD12	2.35	0.52
1:A:201:THR:HG23	1:A:202:PRO:CD	2.40	0.51
1:C:112:PRO:HB2	1:C:115:LYS:HG2	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:25:VAL:HG22	4:D:569:HOH:O	2.10	0.51
1:C:49:LYS:HE2	1:C:185:THR:HG22	1.91	0.51
1:D:264:ARG:HG2	1:D:264:ARG:NH1	2.25	0.51
1:B:125:ILE:HG22	1:D:219:ARG:CG	2.41	0.51
1:D:12:PRO:CG	1:D:202:PRO:HG3	2.37	0.51
1:D:297:ARG:HD3	1:D:326:HIS:ND1	2.25	0.51
1:B:331:LEU:HD11	4:B:564:HOH:O	2.11	0.51
1:D:147:GLU:O	1:D:151:GLU:HG3	2.11	0.51
1:A:137:ASN:HD22	1:A:137:ASN:C	2.13	0.51
1:B:108:ASP:OD1	1:B:110:THR:HG23	2.10	0.51
1:B:219:ARG:HE	1:D:129:ARG:HD3	1.75	0.51
1:D:286:ASN:O	1:D:287:ALA:C	2.53	0.51
1:C:33:ASP:OD1	1:C:232:GLY:HA3	2.11	0.51
1:D:319:GLN:O	1:D:322:GLU:HB2	2.11	0.50
1:A:32:ASP:O	1:A:34:PRO:HD3	2.11	0.50
1:A:195:LEU:O	1:A:201:THR:HB	2.11	0.50
1:C:284:GLU:OE1	1:C:305:PRO:HG3	2.11	0.50
1:D:317:ARG:HG3	1:D:317:ARG:HH11	1.76	0.50
1:B:136:GLU:OE2	1:B:139:HIS:HD2	1.94	0.50
1:B:54:PHE:C	1:B:55:LEU:HG	2.37	0.50
1:B:78:VAL:O	1:B:102:PRO:HA	2.12	0.50
1:D:164:HIS:CD2	4:D:644:HOH:O	2.64	0.50
1:C:137:ASN:O	1:C:138:THR:HB	2.10	0.50
1:B:195:LEU:C	1:B:201:THR:HG22	2.37	0.49
1:C:26:GLY:HA3	1:C:32:ASP:C	2.37	0.49
1:D:132:LEU:HD12	1:D:162:ALA:C	2.36	0.49
1:B:219:ARG:HH21	1:D:129:ARG:HG2	1.76	0.49
1:B:120:ILE:CD1	1:B:159:ARG:HD3	2.43	0.49
1:C:225:GLY:HA3	1:D:60:GLN:HB2	1.93	0.49
1:C:28:ASP:HB3	1:C:228:MET:O	2.12	0.49
1:B:53:ILE:CG2	1:B:207:LEU:HB3	2.42	0.49
1:D:138:THR:HG21	4:D:515:HOH:O	2.12	0.49
1:B:145:PRO:O	1:B:148:TYR:HB3	2.13	0.49
1:B:211:ARG:HB2	1:B:211:ARG:CZ	2.43	0.49
1:B:258:TRP:O	1:B:261:GLU:HB3	2.13	0.49
1:A:279:PHE:CD1	1:A:279:PHE:N	2.81	0.49
1:B:53:ILE:HD11	1:B:221:ARG:HD3	1.94	0.49
1:D:248:ARG:HG3	1:D:248:ARG:HH21	1.77	0.49
1:A:302:ASN:ND2	4:A:568:HOH:O	2.45	0.49
1:D:129:ARG:HG3	1:D:129:ARG:HH11	1.78	0.49
1:D:115:LYS:HD2	1:D:118:MET:CE	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:LEU:HB3	1:A:56:PRO:HD2	1.95	0.48
1:C:234:LEU:HD21	1:D:234:LEU:HD21	1.94	0.48
1:D:239:MET:HE3	4:D:635:HOH:O	2.12	0.48
1:C:137:ASN:ND2	1:C:144:LEU:HB2	2.27	0.48
1:C:244:ASN:C	1:C:245:ASN:HD22	2.22	0.48
1:B:256:THR:HG21	1:B:276:ASN:C	2.39	0.48
1:D:14:ARG:HD3	4:D:571:HOH:O	2.14	0.48
1:D:23:ALA:HB2	1:D:236:ALA:HB2	1.94	0.48
1:D:297:ARG:O	1:D:298:ASN:HB2	2.14	0.48
1:D:213:TYR:OH	4:D:644:HOH:O	2.20	0.48
1:B:292:GLU:O	1:B:293:TYR:C	2.57	0.47
1:A:78:VAL:CG2	1:A:79:GLY:N	2.77	0.47
1:C:110:THR:HB	1:C:148:TYR:CZ	2.49	0.47
1:D:137:ASN:HD22	1:D:138:THR:H	1.60	0.47
1:D:249:LEU:HD22	4:D:551:HOH:O	2.14	0.47
1:A:319:GLN:O	1:A:323:VAL:HG23	2.13	0.47
1:B:54:PHE:HA	1:B:206:LEU:HD23	1.96	0.47
1:B:115:LYS:HB3	1:B:115:LYS:HZ2	1.78	0.47
1:C:36:VAL:HG23	1:C:232:GLY:HA2	1.96	0.47
1:D:138:THR:HG21	1:D:169:ARG:CB	2.43	0.47
1:B:256:THR:HG22	1:B:311:THR:HG21	1.95	0.47
1:A:75:GLU:HB2	1:A:99:GLN:O	2.15	0.47
1:C:169:ARG:NH2	1:C:197:LLP:HE3	2.29	0.47
1:D:56:PRO:HD2	1:D:60:GLN:NE2	2.30	0.47
1:D:262:GLN:HA	1:D:262:GLN:NE2	2.29	0.47
1:C:137:ASN:HD22	1:C:138:THR:N	2.12	0.47
1:D:278:LEU:HD12	1:D:278:LEU:C	2.39	0.47
1:D:69:HIS:CE1	4:D:644:HOH:O	2.68	0.47
1:B:21:MET:O	1:B:21:MET:HG2	2.15	0.46
1:B:256:THR:CG2	1:B:276:ASN:HA	2.42	0.46
1:C:78:VAL:HG22	1:C:79:GLY:N	2.30	0.46
1:D:78:VAL:CG2	1:D:79:GLY:N	2.78	0.46
1:A:234:LEU:HD11	1:B:234:LEU:HD11	1.97	0.46
1:D:236:ALA:HA	1:D:239:MET:CE	2.30	0.46
1:C:2:ILE:CD1	1:C:319:GLN:HB3	2.46	0.46
1:C:196:SER:HA	1:C:201:THR:O	2.14	0.46
1:C:266:ALA:HA	1:C:328:ARG:HE	1.79	0.46
1:D:59:THR:CG2	1:D:88:GLU:HG2	2.26	0.46
1:A:4:LEU:HD21	1:A:315:VAL:HG21	1.98	0.46
1:A:192:THR:HG21	4:A:658:HOH:O	2.14	0.46
1:D:263:LEU:O	1:D:268:ALA:HB3	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:259:MET:HE2	1:A:320:LEU:HD22	1.98	0.46
1:B:180:GLU:HB2	1:B:183:GLU:HG3	1.98	0.46
1:C:1:MET:SD	1:C:300:LEU:HD21	2.56	0.46
1:C:223:MET:HG2	1:D:95:LEU:HD11	1.98	0.46
1:D:138:THR:CG2	4:D:515:HOH:O	2.64	0.46
1:A:121:LYS:HA	1:A:122:PRO:HD3	1.84	0.46
1:D:278:LEU:CD1	1:D:309:LEU:HD12	2.35	0.46
1:D:297:ARG:HD3	1:D:326:HIS:CE1	2.51	0.46
1:B:110:THR:HG21	1:B:145:PRO:HD2	1.97	0.46
1:C:60:GLN:O	1:C:64:VAL:HG23	2.15	0.46
1:C:146:ARG:HG3	1:C:177:TYR:CD1	2.51	0.46
1:A:191:PHE:CE1	1:A:208:VAL:HB	2.51	0.46
1:B:60:GLN:HE22	1:B:226:GLY:CA	2.11	0.46
1:B:110:THR:CG2	1:B:145:PRO:HD2	2.46	0.45
1:B:284:GLU:HG2	4:B:593:HOH:O	2.16	0.45
1:C:279:PHE:HA	1:C:307:VAL:O	2.16	0.45
1:D:271:MET:HE3	1:D:281:ARG:NH2	2.31	0.45
1:C:126:HIS:CE1	4:C:623:HOH:O	2.70	0.45
1:B:191:PHE:HE2	1:B:193:ILE:HD11	1.81	0.45
1:B:191:PHE:CE1	1:B:208:VAL:HB	2.51	0.45
1:A:308:ARG:HD3	4:A:707:HOH:O	2.16	0.45
1:B:186:GLN:C	4:B:548:HOH:O	2.59	0.45
1:C:112:PRO:O	1:C:116:VAL:HG23	2.17	0.45
1:A:155:PHE:CE2	1:A:159:ARG:HD2	2.52	0.45
1:A:138:THR:HG21	4:A:673:HOH:O	2.17	0.45
1:B:121:LYS:HE3	1:C:86:LEU:CD1	2.45	0.45
1:D:47:SER:O	1:D:185:THR:CG2	2.65	0.45
1:A:62:ASN:O	1:A:66:LEU:HG	2.17	0.45
1:D:195:LEU:HD23	1:D:199:LEU:HD12	1.98	0.45
1:C:152:ALA:O	1:C:156:THR:HG23	2.17	0.45
1:B:8:THR:CG2	1:B:197:LLP:HE3	2.47	0.44
1:B:211:ARG:NH1	1:B:211:ARG:CB	2.81	0.44
1:C:191:PHE:CE1	1:C:208:VAL:HB	2.52	0.44
1:A:322:GLU:O	1:A:325:ALA:HB3	2.16	0.44
1:D:115:LYS:HD2	1:D:118:MET:HE1	2.00	0.44
4:A:695:HOH:O	1:B:229:ARG:HG2	2.17	0.44
1:B:261:GLU:O	1:B:265:GLU:HG3	2.18	0.44
1:B:286:ASN:O	1:B:287:ALA:C	2.61	0.44
1:B:86:LEU:HD23	1:B:87:PHE:CE1	2.53	0.44
1:D:259:MET:HE3	1:D:263:LEU:HD21	1.99	0.44
1:A:137:ASN:O	1:A:138:THR:HB	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:195:LEU:C	1:A:201:THR:HB	2.42	0.44
1:C:286:ASN:O	1:C:287:ALA:C	2.61	0.44
1:D:27:ASP:H	1:D:32:ASP:HB2	1.82	0.44
1:C:202:PRO:C	1:C:203:VAL:CG1	2.90	0.44
1:D:258:TRP:CB	1:D:317:ARG:NH1	2.80	0.44
1:A:14:ARG:HB3	4:A:575:HOH:O	2.18	0.44
1:A:35:THR:HG23	1:A:239:MET:CE	2.48	0.44
1:A:59:THR:CG2	1:A:88:GLU:HG2	2.42	0.44
1:A:121:LYS:HE3	1:D:86:LEU:CD1	2.48	0.44
1:B:171:PHE:HB2	4:B:619:HOH:O	2.17	0.44
1:B:256:THR:CG2	1:B:311:THR:OG1	2.66	0.44
1:D:152:ALA:O	1:D:156:THR:HG23	2.18	0.44
1:B:70:CYS:HB2	1:B:98:ILE:HD13	2.00	0.44
1:C:110:THR:HB	1:C:148:TYR:CE1	2.52	0.44
1:C:137:ASN:ND2	1:C:144:LEU:H	2.16	0.44
1:D:240:TYR:CD1	1:D:240:TYR:C	2.95	0.44
1:D:290:LEU:HD23	1:D:290:LEU:C	2.43	0.44
1:A:2:ILE:CD1	1:A:319:GLN:HB3	2.48	0.44
1:B:16:MET:O	1:B:20:MET:HG3	2.18	0.44
1:D:33:ASP:CG	1:D:36:VAL:HG23	2.43	0.44
1:A:167:GLY:H	1:A:192:THR:HG22	1.82	0.43
1:B:201:THR:HG23	1:B:203:VAL:N	2.33	0.43
1:A:137:ASN:C	1:A:137:ASN:ND2	2.76	0.43
1:A:240:TYR:CD1	1:A:240:TYR:C	2.96	0.43
1:B:16:MET:HE1	1:B:200:GLY:C	2.42	0.43
1:B:159:ARG:NH1	1:B:159:ARG:HG2	2.34	0.43
1:C:159:ARG:HE	1:C:159:ARG:HB3	1.65	0.43
1:D:271:MET:HE3	1:D:281:ARG:HH22	1.83	0.43
1:B:124:ASP:HB3	1:B:127:PHE:CD2	2.53	0.43
1:C:70:CYS:HA	1:C:74:GLU:OE1	2.18	0.43
1:C:278:LEU:CD2	1:C:309:LEU:HB2	2.48	0.43
1:C:229:ARG:HB3	1:C:230:GLN:HE22	1.81	0.43
1:A:259:MET:HE2	1:A:320:LEU:CD2	2.48	0.43
1:B:27:ASP:H	1:B:32:ASP:HB2	1.83	0.43
1:B:170:ILE:O	1:B:173:ALA:HB3	2.18	0.43
1:B:191:PHE:CE2	1:B:193:ILE:HD11	2.54	0.43
1:C:53:ILE:HG23	1:C:55:LEU:HD11	2.00	0.43
1:C:167:GLY:O	1:C:168:ALA:C	2.62	0.43
1:D:286:ASN:HD22	1:D:286:ASN:N	2.15	0.43
1:A:168:ALA:HA	1:A:194:CYS:HB2	2.00	0.43
1:C:121:LYS:HA	1:C:122:PRO:HD3	1.90	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:138:THR:HG22	4:D:565:HOH:O	2.18	0.43
1:C:248:ARG:NH2	1:C:252:ASP:OD2	2.49	0.43
1:A:110:THR:HG22	1:A:144:LEU:HG	2.01	0.43
1:C:2:ILE:HD13	1:C:319:GLN:HB3	2.00	0.43
1:C:248:ARG:NE	4:C:539:HOH:O	2.40	0.43
1:D:139:HIS:HE1	4:D:589:HOH:O	2.01	0.43
1:D:216:ARG:HD2	4:D:577:HOH:O	2.18	0.43
1:A:70:CYS:HB2	1:A:98:ILE:HD13	2.00	0.42
1:A:120:ILE:CD1	1:A:159:ARG:HD3	2.48	0.42
1:A:196:SER:HA	1:A:201:THR:O	2.19	0.42
1:B:42:TYR:O	1:B:45:GLU:HB3	2.19	0.42
1:B:138:THR:CG2	4:B:504:HOH:O	2.67	0.42
1:C:256:THR:HG23	1:C:278:LEU:HD22	2.01	0.42
1:A:304:SER:OG	1:A:305:PRO:HD2	2.20	0.42
1:B:276:ASN:HB2	1:B:311:THR:OG1	2.19	0.42
1:C:279:PHE:N	1:C:279:PHE:CD1	2.87	0.42
1:C:316:SER:OG	1:C:319:GLN:HG3	2.20	0.42
1:A:248:ARG:HG3	1:A:248:ARG:HH11	1.85	0.42
1:B:219:ARG:HG3	1:B:219:ARG:HH11	1.85	0.42
1:C:137:ASN:HD22	1:C:138:THR:H	1.67	0.42
1:C:202:PRO:O	1:C:203:VAL:HG12	2.20	0.42
1:D:259:MET:HE3	1:D:263:LEU:HD11	2.01	0.42
1:A:284:GLU:OE2	1:A:284:GLU:O	2.38	0.42
1:B:20:MET:O	1:B:233:ILE:HG13	2.20	0.42
1:D:35:THR:HG23	1:D:239:MET:CE	2.50	0.42
1:D:193:ILE:O	1:D:205:SER:HA	2.20	0.42
1:D:301:ILE:C	1:D:301:ILE:CD1	2.90	0.42
1:A:319:GLN:O	1:A:322:GLU:HB2	2.21	0.41
1:C:50:GLU:HG3	4:C:530:HOH:O	2.19	0.41
1:C:147:GLU:CD	1:C:147:GLU:H	2.27	0.41
1:C:183:GLU:HA	1:C:186:GLN:HE21	1.84	0.41
1:B:138:THR:CA	1:B:142:LYS:O	2.68	0.41
1:A:236:ALA:HA	1:A:239:MET:HE2	2.01	0.41
4:A:680:HOH:O	1:B:222:LYS:NZ	2.35	0.41
1:C:119:LYS:HD3	1:C:119:LYS:HA	1.85	0.41
1:C:255:ASN:HD22	1:C:317:ARG:HH21	1.69	0.41
1:B:219:ARG:HH12	1:D:125:ILE:HA	1.82	0.41
1:C:195:LEU:C	1:C:201:THR:CG2	2.93	0.41
1:C:278:LEU:HD21	1:C:309:LEU:HB2	2.03	0.41
1:D:132:LEU:CD1	1:D:162:ALA:HB3	2.51	0.41
1:C:131:LYS:NZ	4:C:651:HOH:O	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:240:TYR:CD1	1:C:240:TYR:C	2.98	0.41
1:C:283:GLY:O	1:C:287:ALA:HB2	2.20	0.41
1:A:262:GLN:OE1	1:A:262:GLN:HA	2.21	0.41
1:C:175:VAL:HG21	1:C:249:LEU:HB2	2.02	0.41
1:C:245:ASN:N	1:C:245:ASN:ND2	2.68	0.41
1:D:269:ASP:OD2	1:D:281:ARG:NH1	2.50	0.41
1:D:278:LEU:HD12	1:D:278:LEU:O	2.20	0.41
1:D:279:PHE:CD1	1:D:279:PHE:N	2.89	0.41
1:A:201:THR:HA	1:A:202:PRO:HD3	1.91	0.41
1:A:233:ILE:HG23	1:A:234:LEU:N	2.35	0.41
1:B:271:MET:O	1:B:272:ARG:HB3	2.21	0.41
1:C:1:MET:HB2	1:C:298:ASN:O	2.21	0.41
1:C:5:ARG:HG2	1:C:302:ASN:ND2	2.35	0.41
1:A:197:LLP:H4'1	3:A:402[A]:SER:N	2.36	0.41
1:A:199:LEU:HD23	1:A:199:LEU:HA	1.85	0.41
1:A:203:VAL:HG22	1:A:234:LEU:HD13	2.00	0.41
1:C:23:ALA:HA	1:C:24:PRO:HD3	1.74	0.41
1:B:115:LYS:NZ	1:B:115:LYS:CB	2.83	0.41
1:B:154:GLU:HB2	4:B:660:HOH:O	2.21	0.41
1:B:290:LEU:HA	1:B:330:PHE:CE1	2.56	0.41
1:B:297:ARG:HD3	1:B:326:HIS:CD2	2.55	0.41
1:C:71:GLU:HG2	4:C:651:HOH:O	2.20	0.41
1:D:33:ASP:OD1	1:D:36:VAL:HG23	2.21	0.41
1:D:203:VAL:CG1	1:D:204:GLY:N	2.83	0.41
1:B:146:ARG:O	1:B:149:LEU:HB2	2.21	0.40
1:C:145:PRO:O	1:C:148:TYR:HB3	2.21	0.40
1:C:166:ASP:HA	1:C:192:THR:HG22	2.02	0.40
1:C:277:MET:SD	1:C:308:ARG:HD2	2.61	0.40
1:C:318:ALA:O	1:C:321:ALA:HB3	2.21	0.40
1:A:139:HIS:HE1	4:A:601:HOH:O	2.03	0.40
1:B:75:GLU:HA	1:B:99:GLN:O	2.22	0.40
1:B:196:SER:HA	1:B:201:THR:O	2.21	0.40
1:C:75:GLU:HA	1:C:99:GLN:O	2.21	0.40
1:D:165:VAL:O	1:D:191:PHE:HA	2.21	0.40
1:A:245:ASN:HB3	4:A:512:HOH:O	2.20	0.40
1:C:167:GLY:N	1:C:192:THR:CG2	2.77	0.40
1:C:263:LEU:HD23	1:C:263:LEU:HA	1.84	0.40
1:D:47:SER:CB	1:D:185:THR:HG21	2.51	0.40
1:B:90:GLY:O	1:B:94:VAL:HG23	2.21	0.40
1:C:78:VAL:CG2	1:C:79:GLY:N	2.84	0.40
1:C:304:SER:O	1:C:305:PRO:C	2.64	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:313:LEU:C	1:D:313:LEU:HD12	2.46	0.40
1:C:201:THR:HG23	1:C:203:VAL:H	1.87	0.40
1:D:75:GLU:OE1	1:D:121:LYS:NZ	2.54	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	329/333 (99%)	318 (97%)	10 (3%)	1 (0%)	36	36
1	B	329/333 (99%)	310 (94%)	17 (5%)	2 (1%)	21	18
1	C	329/333 (99%)	316 (96%)	11 (3%)	2 (1%)	21	18
1	D	329/333 (99%)	313 (95%)	15 (5%)	1 (0%)	36	36
All	All	1316/1332 (99%)	1257 (96%)	53 (4%)	6 (0%)	24	22

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	138	THR
1	A	138	THR
1	C	138	THR
1	D	138	THR
1	B	293	TYR
1	C	229	ARG

5.3.2 Protein sidechains [i](#)

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	259/260 (100%)	252 (97%)	7 (3%)	39	45
1	B	259/260 (100%)	247 (95%)	12 (5%)	24	25
1	C	259/260 (100%)	245 (95%)	14 (5%)	20	18
1	D	259/260 (100%)	253 (98%)	6 (2%)	44	51
All	All	1036/1040 (100%)	997 (96%)	39 (4%)	28	32

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	ILE
1	A	114	ASP
1	A	115	LYS
1	A	138	THR
1	A	147	GLU
1	A	201	THR
1	A	203	VAL
1	B	53	ILE
1	B	115	LYS
1	B	138	THR
1	B	144	LEU
1	B	147	GLU
1	B	185	THR
1	B	201	THR
1	B	203	VAL
1	B	205	SER
1	B	219	ARG
1	B	256	THR
1	B	315	VAL
1	C	53	ILE
1	C	99	GLN
1	C	138	THR
1	C	144	LEU
1	C	147	GLU
1	C	157	ARG
1	C	201	THR
1	C	203	VAL
1	C	230	GLN
1	C	233	ILE
1	C	239	MET

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Mol	Chain	Res	Type
1	C	248	ARG
1	C	273	GLN
1	C	297	ARG
1	D	45	GLU
1	D	67	LEU
1	D	69	HIS
1	D	138	THR
1	D	147	GLU
1	D	273	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (50) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	60	GLN
1	A	137	ASN
1	A	139	HIS
1	A	140	ASN
1	A	164	HIS
1	A	210	ASN
1	A	230	GLN
1	A	245	ASN
1	A	302	ASN
1	B	60	GLN
1	B	139	HIS
1	B	172	ASN
1	B	210	ASN
1	B	230	GLN
1	B	245	ASN
1	B	273	GLN
1	B	302	ASN
1	B	326	HIS
1	C	60	GLN
1	C	99	GLN
1	C	137	ASN
1	C	140	ASN
1	C	164	HIS
1	C	172	ASN
1	C	186	GLN
1	C	210	ASN
1	C	230	GLN
1	C	245	ASN

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Mol	Chain	Res	Type
1	C	250	GLN
1	C	255	ASN
1	C	262	GLN
1	C	273	GLN
1	C	298	ASN
1	C	302	ASN
1	D	37	ASN
1	D	60	GLN
1	D	69	HIS
1	D	137	ASN
1	D	139	HIS
1	D	164	HIS
1	D	172	ASN
1	D	210	ASN
1	D	230	GLN
1	D	244	ASN
1	D	245	ASN
1	D	250	GLN
1	D	262	GLN
1	D	286	ASN
1	D	302	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	LLP	D	197	1	23,24,25	3.72	8 (34%)	25,32,34	2.74	11 (44%)
1	LLP	B	197	1	23,24,25	3.72	9 (39%)	25,32,34	2.69	10 (40%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	LLP	C	197	1	23,24,25	3.86	9 (39%)	25,32,34	2.54	10 (40%)
1	LLP	A	197	1	23,24,25	3.86	7 (30%)	25,32,34	2.67	10 (40%)

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
1	LLP	D	197	1	-	7/16/17/19	0/1/1/1
1	LLP	B	197	1	-	9/16/17/19	0/1/1/1
1	LLP	C	197	1	-	10/16/17/19	0/1/1/1
1	LLP	A	197	1	-	4/16/17/19	0/1/1/1

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	197	LLP	C4-C5	13.62	1.60	1.42
1	A	197	LLP	C4-C5	13.49	1.60	1.42
1	D	197	LLP	C4-C5	13.14	1.60	1.42
1	B	197	LLP	C4-C5	12.75	1.59	1.42
1	A	197	LLP	C3-C2	8.04	1.49	1.41
1	C	197	LLP	C3-C2	7.99	1.49	1.41
1	B	197	LLP	C3-C2	7.88	1.49	1.41
1	D	197	LLP	C3-C2	7.61	1.48	1.41
1	C	197	LLP	C2-N1	5.27	1.43	1.33
1	A	197	LLP	C2-N1	5.21	1.43	1.33
1	D	197	LLP	C2-N1	4.96	1.42	1.33
1	B	197	LLP	C2-N1	4.94	1.42	1.33
1	C	197	LLP	C6-N1	4.42	1.43	1.34
1	B	197	LLP	C6-N1	3.99	1.42	1.34
1	A	197	LLP	C4'-NZ	3.92	1.40	1.27
1	D	197	LLP	C6-N1	3.88	1.42	1.34
1	A	197	LLP	C6-N1	3.74	1.42	1.34
1	D	197	LLP	C4'-NZ	3.56	1.39	1.27
1	B	197	LLP	C4-C3	3.39	1.46	1.41
1	C	197	LLP	C4'-NZ	3.34	1.38	1.27
1	D	197	LLP	C4-C3	3.08	1.46	1.41
1	A	197	LLP	C4-C3	3.02	1.46	1.41
1	B	197	LLP	C4'-NZ	2.96	1.37	1.27
1	B	197	LLP	C6-C5	2.58	1.42	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	197	LLP	P-OP3	-2.49	1.45	1.54
1	B	197	LLP	C5'-C5	2.48	1.57	1.50
1	C	197	LLP	C5'-C5	2.29	1.56	1.50
1	D	197	LLP	P-OP3	-2.19	1.46	1.54
1	C	197	LLP	C6-C5	2.17	1.42	1.37
1	C	197	LLP	C4-C3	2.15	1.44	1.41
1	B	197	LLP	P-OP4	2.07	1.66	1.60
1	A	197	LLP	OP4-C5'	-2.04	1.37	1.44
1	D	197	LLP	C6-C5	2.03	1.41	1.37

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	197	LLP	C2'-C2-C3	8.56	130.81	120.80
1	D	197	LLP	C2'-C2-C3	8.15	130.34	120.80
1	B	197	LLP	C2'-C2-C3	8.01	130.17	120.80
1	C	197	LLP	C2'-C2-C3	7.91	130.05	120.80
1	B	197	LLP	OP4-C5'-C5	5.23	119.17	109.36
1	D	197	LLP	OP4-C5'-C5	4.72	118.20	109.36
1	D	197	LLP	C6-N1-C2	4.38	127.14	119.20
1	A	197	LLP	C6-N1-C2	4.33	127.05	119.20
1	B	197	LLP	C6-N1-C2	4.17	126.77	119.20
1	C	197	LLP	C6-N1-C2	4.14	126.70	119.20
1	A	197	LLP	OP4-C5'-C5	3.91	116.69	109.36
1	A	197	LLP	C5-C4-C4'	3.87	127.44	121.47
1	C	197	LLP	C5-C4-C4'	3.73	127.23	121.47
1	D	197	LLP	C5-C4-C4'	3.73	127.23	121.47
1	C	197	LLP	OP4-C5'-C5	3.69	116.27	109.36
1	B	197	LLP	C5-C4-C4'	3.46	126.80	121.47
1	A	197	LLP	C3-C2-N1	-3.33	116.76	120.96
1	B	197	LLP	C3-C2-N1	-3.28	116.82	120.96
1	C	197	LLP	C5-C6-N1	-3.23	118.58	123.83
1	D	197	LLP	C3-C2-N1	-3.23	116.89	120.96
1	C	197	LLP	C3-C4-C4'	-3.21	114.60	120.40
1	D	197	LLP	C5-C6-N1	-3.08	118.82	123.83
1	C	197	LLP	C3-C2-N1	-3.02	117.15	120.96
1	A	197	LLP	C3-C4-C4'	-2.97	115.04	120.40
1	D	197	LLP	C3-C4-C4'	-2.95	115.08	120.40
1	A	197	LLP	C5-C6-N1	-2.92	119.08	123.83
1	A	197	LLP	C2'-C2-N1	-2.77	112.42	117.64
1	B	197	LLP	C3-C4-C4'	-2.75	115.44	120.40
1	B	197	LLP	C5-C6-N1	-2.73	119.39	123.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	197	LLP	O3-C3-C4	2.66	126.61	119.44
1	D	197	LLP	O3-C3-C4	2.65	126.58	119.44
1	D	197	LLP	C2'-C2-N1	-2.58	112.77	117.64
1	A	197	LLP	O3-C3-C4	2.57	126.37	119.44
1	C	197	LLP	C2'-C2-N1	-2.57	112.80	117.64
1	B	197	LLP	O3-C3-C2	-2.49	112.42	117.58
1	B	197	LLP	C2'-C2-N1	-2.46	113.02	117.64
1	C	197	LLP	O3-C3-C4	2.43	125.98	119.44
1	D	197	LLP	C4-C4'-NZ	-2.37	113.12	124.04
1	D	197	LLP	O3-C3-C2	-2.33	112.75	117.58
1	A	197	LLP	O3-C3-C2	-2.17	113.08	117.58
1	C	197	LLP	O3-C3-C2	-2.03	113.36	117.58

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	197	LLP	C4-C4'-NZ-CE
1	A	197	LLP	O-C-CA-CB
1	B	197	LLP	C4-C4'-NZ-CE
1	B	197	LLP	C4-C5-C5'-OP4
1	B	197	LLP	C6-C5-C5'-OP4
1	B	197	LLP	C5'-OP4-P-OP2
1	B	197	LLP	C5'-OP4-P-OP3
1	B	197	LLP	O-C-CA-CB
1	C	197	LLP	C4-C4'-NZ-CE
1	C	197	LLP	C4-C5-C5'-OP4
1	C	197	LLP	C6-C5-C5'-OP4
1	C	197	LLP	C5'-OP4-P-OP1
1	C	197	LLP	C5'-OP4-P-OP2
1	C	197	LLP	C5'-OP4-P-OP3
1	C	197	LLP	O-C-CA-CB
1	D	197	LLP	C4-C4'-NZ-CE
1	D	197	LLP	C4-C5-C5'-OP4
1	D	197	LLP	C6-C5-C5'-OP4
1	D	197	LLP	O-C-CA-CB
1	A	197	LLP	C3-C4-C4'-NZ
1	C	197	LLP	C3-C4-C4'-NZ
1	B	197	LLP	C5'-OP4-P-OP1
1	C	197	LLP	CG-CD-CE-NZ
1	D	197	LLP	CD-CE-NZ-C4'
1	A	197	LLP	CG-CD-CE-NZ

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Mol	Chain	Res	Type	Atoms
1	B	197	LLP	C3-C4-C4'-NZ
1	D	197	LLP	C3-C4-C4'-NZ
1	C	197	LLP	CD-CE-NZ-C4'
1	D	197	LLP	C5'-OP4-P-OP2
1	B	197	LLP	CD-CE-NZ-C4'

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	197	LLP	2	0
1	C	197	LLP	4	0
1	A	197	LLP	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 14 ligands modelled in this entry, 6 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	SER	D	403[A]	-	4,6,6	1.12	0	2,7,7	0.47	0
3	SER	A	402[A]	-	4,6,6	1.35	0	2,7,7	0.59	0
3	SER	A	402[B]	-	4,6,6	1.44	0	2,7,7	0.27	0
3	SER	C	402[B]	-	4,6,6	1.25	0	2,7,7	0.26	0
3	SER	D	403[B]	-	4,6,6	1.34	0	2,7,7	0.49	0
3	SER	B	403[A]	-	4,6,6	1.24	0	2,7,7	0.27	0
3	SER	B	403[B]	-	4,6,6	1.19	0	2,7,7	0.24	0
3	SER	C	402[A]	-	4,6,6	1.14	0	2,7,7	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	SER	D	403[A]	-	-	3/6/6/6	-
3	SER	A	402[A]	-	-	6/6/6/6	-
3	SER	A	402[B]	-	-	2/6/6/6	-
3	SER	C	402[B]	-	-	2/6/6/6	-
3	SER	D	403[B]	-	-	6/6/6/6	-
3	SER	B	403[A]	-	-	2/6/6/6	-
3	SER	B	403[B]	-	-	2/6/6/6	-
3	SER	C	402[A]	-	-	0/6/6/6	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	402[A]	SER	O-C-CA-CB
3	A	402[A]	SER	OXT-C-CA-CB
3	A	402[A]	SER	N-CA-CB-OG
3	A	402[A]	SER	C-CA-CB-OG
3	B	403[A]	SER	C-CA-CB-OG
3	D	403[A]	SER	OXT-C-CA-CB
3	D	403[B]	SER	O-C-CA-CB
3	D	403[B]	SER	OXT-C-CA-CB
3	D	403[B]	SER	N-CA-CB-OG
3	D	403[B]	SER	C-CA-CB-OG
3	D	403[B]	SER	OXT-C-CA-N
3	A	402[A]	SER	OXT-C-CA-N
3	A	402[B]	SER	OXT-C-CA-N
3	B	403[B]	SER	OXT-C-CA-N
3	C	402[B]	SER	OXT-C-CA-N
3	A	402[B]	SER	O-C-CA-N
3	C	402[B]	SER	O-C-CA-N
3	B	403[A]	SER	N-CA-CB-OG
3	D	403[A]	SER	N-CA-CB-OG
3	D	403[A]	SER	C-CA-CB-OG
3	A	402[A]	SER	O-C-CA-N

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Mol	Chain	Res	Type	Atoms
3	B	403[B]	SER	O-C-CA-N
3	D	403[B]	SER	O-C-CA-N

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	402[A]	SER	1	0
3	B	403[A]	SER	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	331/333 (99%)	-1.68	0 100 100	12, 23, 37, 52	0
1	B	331/333 (99%)	-1.64	0 100 100	14, 26, 46, 59	0
1	C	331/333 (99%)	-1.63	0 100 100	12, 27, 42, 58	0
1	D	331/333 (99%)	-1.68	0 100 100	12, 24, 39, 58	0
All	All	1324/1332 (99%)	-1.66	0 100 100	12, 25, 42, 59	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	LLP	A	197	24/25	1.00	0.02	16,21,33,34	0
1	LLP	B	197	24/25	1.00	0.02	18,30,35,36	0
1	LLP	C	197	24/25	1.00	0.02	16,25,33,38	0
1	LLP	D	197	24/25	1.00	0.02	14,27,31,36	0

6.3 Carbohydrates [i](#)

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(Å ²)	Q<0.9
3	SER	B	403[A]	7/7	0.98	0.04	30,39,44,44	7
3	SER	B	403[B]	7/7	0.98	0.04	25,32,37,38	7
3	SER	A	402[A]	7/7	0.99	0.04	16,24,32,33	7
3	SER	A	402[B]	7/7	0.99	0.04	14,18,23,25	7
2	NA	B	402	1/1	0.99	0.02	19,19,19,19	0
2	NA	D	402	1/1	0.99	0.03	28,28,28,28	0
3	SER	C	402[A]	7/7	0.99	0.05	37,42,46,46	7
3	SER	C	402[B]	7/7	0.99	0.05	29,33,36,37	7
3	SER	D	403[A]	7/7	0.99	0.03	22,32,35,37	7
3	SER	D	403[B]	7/7	0.99	0.03	19,24,30,31	7
2	NA	A	401	1/1	1.00	0.01	23,23,23,23	0
2	NA	C	401	1/1	1.00	0.01	23,23,23,23	0
2	NA	D	401	1/1	1.00	0.05	22,22,22,22	0
2	NA	B	401	1/1	1.00	0.04	14,14,14,14	0

6.5 Other polymers

There are no such residues in this entry.