



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:20 PM UTC

PDB ID : 4ECA / pdb_00004eca
Title : ASPARAGINASE FROM E. COLI, MUTANT T89V WITH COVALENTLY BOUND ASPARTATE
Authors : Palm, G.J.; Lubkowski, J.; Wlodawer, A.
Deposited on : 1997-02-21
Resolution : 2.20 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

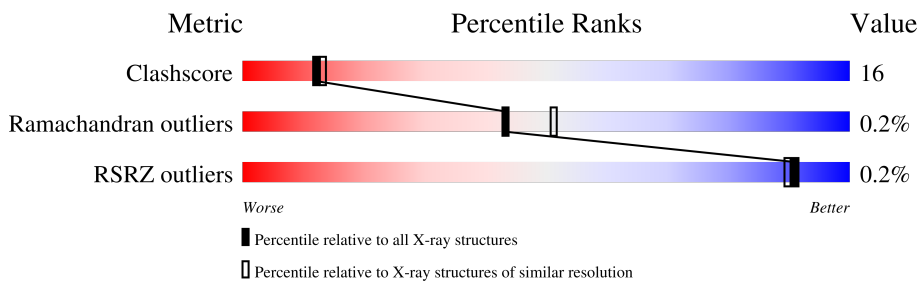
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	326	
1	B	326	
1	C	326	
1	D	326	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10476 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 L-ASPARAGINE AMIDOHYDROLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	326	2439	1522	416	493	8	0	0	0
1	B	326	2439	1522	416	493	8	0	0	0
1	C	326	2439	1522	416	493	8	0	0	0
1	D	326	2439	1522	416	493	8	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	12	AEI	THR	modified residue	UNP P00805
A	89	VAL	THR	engineered mutation	UNP P00805
B	12	AEI	THR	modified residue	UNP P00805
B	89	VAL	THR	engineered mutation	UNP P00805
C	12	AEI	THR	modified residue	UNP P00805
C	89	VAL	THR	engineered mutation	UNP P00805
D	12	AEI	THR	modified residue	UNP P00805
D	89	VAL	THR	engineered mutation	UNP P00805

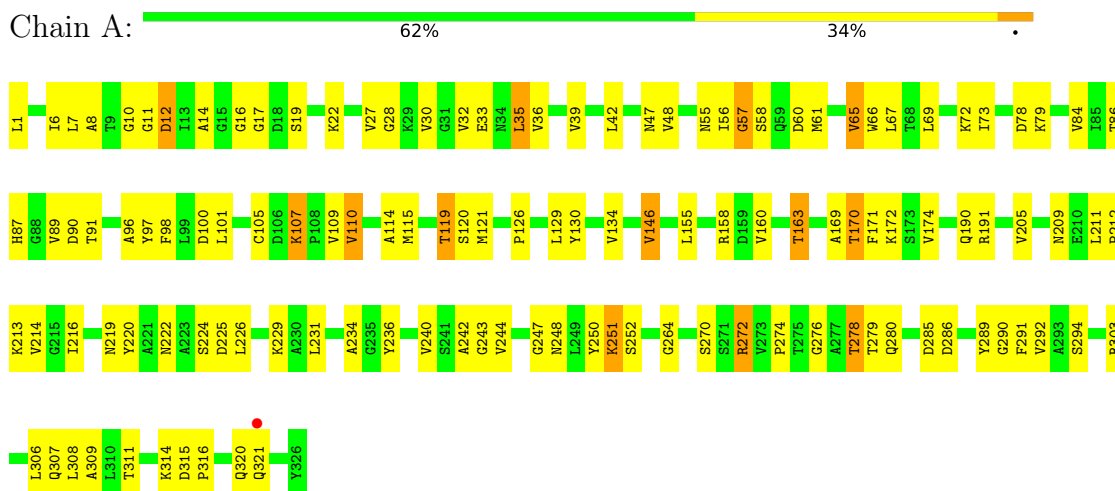
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	191	Total 191	O 191	0	0
2	B	172	Total 172	O 172	0	0
2	C	187	Total 187	O 187	0	0
2	D	170	Total 170	O 170	0	0

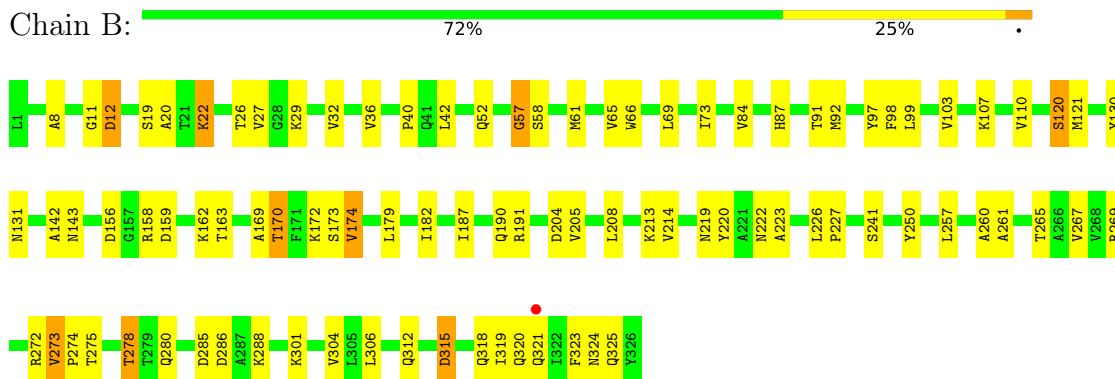
3 Residue-property plots

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.

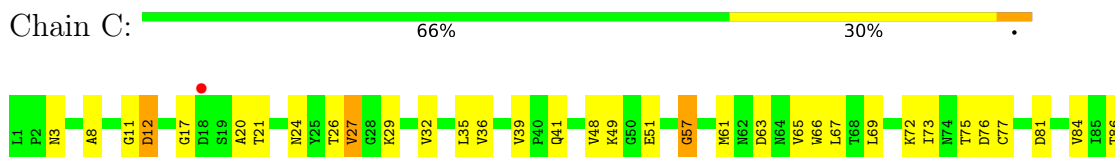
• Molecule 1: L-ASPARAGINE AMIDOHYDROLASE

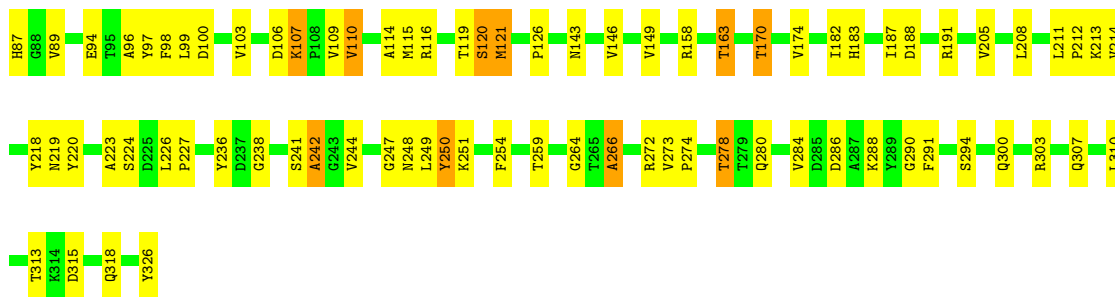


• Molecule 1: L-ASPARAGINE AMIDOHYDROLASE

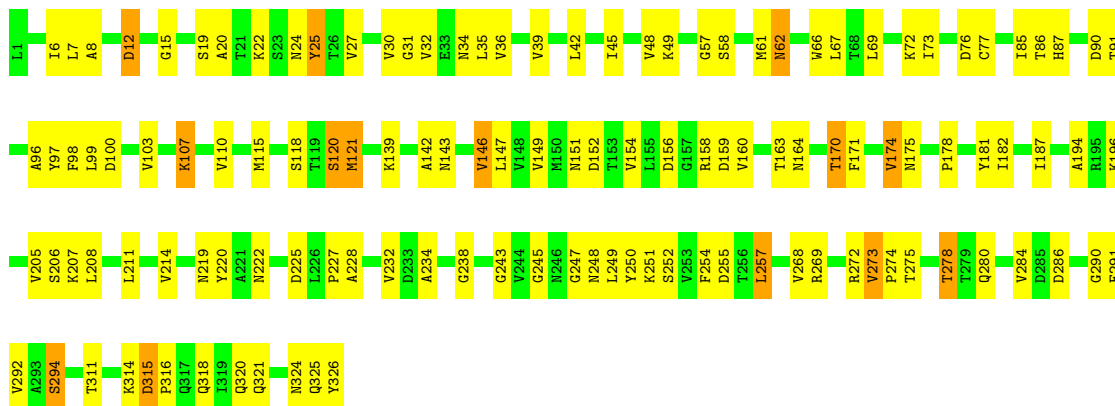


• Molecule 1: L-ASPARAGINE AMIDOHYDROLASE





● Molecule 1: L-ASPARAGINE AMIDOHYDROLASE



4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	95.00Å 126.20Å 155.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.20 10.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	62.4 (10.00-2.20) 63.1 (10.00-2.20)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.11 (at 2.00Å)	Xtrriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.181 , (Not available) 0.176 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	17.5	Xtrriage
Anisotropy	0.448	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 63.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10476	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 46.65 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1105e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality i

5.1 Standard geometry i

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

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	0.90	2/2460 (0.1%)	1.19	24/3348 (0.7%)
1	B	0.90	2/2460 (0.1%)	1.20	13/3348 (0.4%)
1	C	0.86	1/2460 (0.0%)	1.24	27/3348 (0.8%)
1	D	0.89	3/2460 (0.1%)	1.21	23/3348 (0.7%)
All	All	0.89	8/9840 (0.1%)	1.21	87/13392 (0.6%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	92	MET	SD-CE	-6.58	1.63	1.79
1	B	121	MET	SD-CE	-6.58	1.63	1.79
1	A	121	MET	SD-CE	-6.34	1.63	1.79
1	A	146	VAL	CA-CB	-5.79	1.47	1.54
1	D	146	VAL	CA-CB	-5.53	1.48	1.54
1	C	121	MET	SD-CE	-5.25	1.66	1.79
1	D	121	MET	SD-CE	-5.22	1.66	1.79
1	D	45	ILE	CA-CB	5.13	1.61	1.54

All (87) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	170	THR	N-CA-C	8.37	120.40	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	174	VAL	N-CA-C	8.32	123.87	111.89
1	C	57	GLY	N-CA-C	-8.12	98.89	111.64
1	A	107	LYS	N-CA-C	-8.02	98.54	110.24
1	C	280	GLN	N-CA-C	-7.99	98.64	110.46
1	C	264	GLY	N-CA-C	7.83	121.96	113.58
1	D	274	PRO	N-CA-C	7.77	123.94	113.84
1	C	170	THR	N-CA-C	7.61	120.53	111.33
1	C	273	VAL	CA-C-N	7.57	127.45	119.05
1	C	273	VAL	C-N-CA	7.57	127.45	119.05
1	B	280	GLN	N-CA-C	-7.55	99.29	110.46
1	D	120	SER	N-CA-C	7.44	120.59	110.55
1	B	278	THR	N-CA-C	-7.32	96.62	108.41
1	A	280	GLN	N-CA-C	-7.27	99.76	110.52
1	A	121	MET	N-CA-C	-7.17	99.67	110.28
1	A	309	ALA	N-CA-C	-7.16	103.47	111.28
1	D	278	THR	N-CA-C	-7.03	97.09	108.41
1	C	48	VAL	N-CA-C	6.98	118.71	109.21
1	C	84	VAL	N-CA-C	-6.91	98.09	108.17
1	A	174	VAL	N-CA-C	6.88	121.80	111.89
1	D	62	ASN	N-CA-C	6.84	118.54	108.14
1	D	174	VAL	N-CA-C	6.83	121.72	111.89
1	C	294	SER	N-CA-C	6.81	121.58	113.28
1	B	120	SER	N-CA-C	6.74	119.65	110.55
1	A	274	PRO	N-CA-C	6.68	122.81	113.53
1	A	47	ASN	N-CA-C	-6.67	95.33	107.75
1	D	315	ASP	CA-C-N	6.64	128.14	119.84
1	D	315	ASP	C-N-CA	6.64	128.14	119.84
1	A	272	ARG	N-CA-C	-6.59	104.50	112.54
1	C	27	VAL	N-CA-C	6.53	118.01	109.58
1	C	242	ALA	N-CA-C	-6.53	97.20	108.56
1	C	163	THR	N-CA-C	6.39	120.28	112.23
1	A	311	THR	N-CA-C	-6.28	105.45	113.23
1	B	107	LYS	N-CA-C	-6.22	101.16	110.24
1	D	273	VAL	CA-C-N	6.12	125.80	119.56
1	D	273	VAL	C-N-CA	6.12	125.80	119.56
1	A	120	SER	N-CA-C	6.08	118.15	110.24
1	D	57	GLY	N-CA-C	-6.04	100.44	111.25
1	B	57	GLY	N-CA-C	-5.97	102.26	111.64
1	C	107	LYS	N-CA-C	-5.96	101.45	110.39
1	A	57	GLY	N-CA-C	-5.95	101.09	111.62
1	C	119	THR	N-CA-C	5.85	118.61	111.82
1	C	120	SER	N-CA-C	5.80	118.03	110.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	174	VAL	N-CA-C	5.80	120.25	111.89
1	D	248	ASN	N-CA-C	5.75	118.61	110.50
1	C	259	THR	N-CA-C	-5.75	104.94	111.14
1	D	49	LYS	N-CA-C	-5.74	99.06	108.76
1	D	280	GLN	N-CA-C	-5.69	101.60	110.42
1	B	274	PRO	N-CA-C	5.69	122.06	113.81
1	D	294	SER	N-CA-C	5.69	119.48	112.54
1	D	25	TYR	N-CA-C	5.65	116.82	107.73
1	D	160	VAL	N-CA-C	5.64	116.98	109.37
1	C	315	ASP	CA-C-N	5.64	125.75	119.32
1	C	315	ASP	C-N-CA	5.64	125.75	119.32
1	B	22	LYS	N-CA-C	-5.61	99.75	108.90
1	B	315	ASP	CA-C-N	5.61	125.07	119.24
1	B	315	ASP	C-N-CA	5.61	125.07	119.24
1	D	170	THR	N-CA-C	5.61	118.11	111.33
1	B	273	VAL	CA-C-N	5.58	125.94	119.47
1	B	273	VAL	C-N-CA	5.58	125.94	119.47
1	A	278	THR	N-CA-C	-5.57	98.65	108.23
1	C	273	VAL	N-CA-C	-5.57	103.49	108.95
1	A	119	THR	N-CA-C	5.55	117.41	111.36
1	C	250	TYR	N-CA-C	-5.54	102.08	110.28
1	A	48	VAL	N-CA-C	5.49	116.48	108.58
1	A	35	LEU	N-CA-C	-5.47	105.40	111.36
1	D	107	LYS	N-CA-C	-5.46	102.20	110.39
1	C	266	ALA	N-CA-C	-5.44	101.61	110.20
1	D	90	ASP	N-CA-C	5.38	117.56	111.11
1	A	110	VAL	N-CA-C	5.37	115.62	108.11
1	D	284	VAL	N-CA-C	5.37	116.03	108.36
1	C	284	VAL	N-CA-C	5.33	116.65	108.71
1	A	279	THR	N-CA-C	5.29	117.69	110.55
1	A	242	ALA	N-CA-C	-5.26	98.10	107.98
1	C	278	THR	N-CA-C	-5.25	98.75	107.99
1	D	257	LEU	N-CA-C	5.24	117.07	111.36
1	D	311	THR	N-CA-C	-5.20	106.78	113.23
1	A	65	VAL	N-CA-C	-5.19	104.61	111.09
1	C	41	GLN	N-CA-C	-5.16	106.49	112.89
1	A	209	ASN	N-CA-C	-5.15	106.40	113.30
1	D	171	PHE	N-CA-C	5.13	116.62	108.67
1	A	163	THR	N-CA-CB	-5.12	102.64	110.33
1	B	170	THR	N-CA-C	5.12	117.60	111.71
1	C	110	VAL	N-CA-C	5.12	115.95	108.53
1	A	321	GLN	N-CA-C	-5.07	105.84	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	264	GLY	N-CA-C	5.06	121.79	115.21
1	C	121	MET	N-CA-C	-5.06	102.21	110.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	326	TYR	Sidechain

5.2 Too-close contacts [i](#)

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	2439	0	2429	84	0
1	B	2439	0	2429	73	0
1	C	2439	0	2429	81	0
1	D	2439	0	2429	82	0
2	A	191	0	0	5	0
2	B	172	0	0	12	0
2	C	187	0	0	6	0
2	D	170	0	0	3	0
All	All	10476	0	9716	303	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 16.

All (303) 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:C:12:AEI:HZ	1:C:27:VAL:HG21	1.53	0.90
1:D:73:ILE:HD11	1:D:85:ILE:HD11	1.54	0.87
1:C:27:VAL:HG11	1:C:57:GLY:HA2	1.62	0.81
1:C:163:THR:HB	1:C:170:THR:O	1.82	0.80
2:B:421:HOH:O	1:C:183:HIS:HD2	1.65	0.80
1:D:321:GLN:O	1:D:325:GLN:HG2	1.82	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:26:THR:HB	1:B:29:LYS:HE3	1.66	0.77
1:C:205:VAL:HA	1:C:208:LEU:HD12	1.64	0.77
1:A:17:GLY:H	1:A:119:THR:HG21	1.50	0.77
1:D:58:SER:HB2	1:D:91:THR:OG1	1.85	0.76
1:B:227:PRO:HB3	1:D:227:PRO:HB3	1.68	0.75
1:D:61:MET:HE3	1:D:87:HIS:NE2	2.02	0.75
1:A:19:SER:HB3	1:A:22:LYS:HB2	1.70	0.72
1:A:163:THR:HB	1:A:170:THR:O	1.90	0.71
1:C:20:ALA:HA	1:C:120:SER:HA	1.71	0.71
1:D:163:THR:HB	1:D:170:THR:O	1.90	0.71
1:D:219:ASN:ND2	1:D:250:TYR:H	1.89	0.70
1:A:32:VAL:O	1:A:36:VAL:HG23	1.93	0.69
1:A:97:TYR:CE1	1:A:158:ARG:HG3	2.28	0.69
1:D:107:LYS:HE2	2:D:477:HOH:O	1.94	0.67
1:A:61:MET:HE3	1:A:87:HIS:NE2	2.10	0.67
1:D:19:SER:HB3	1:D:22:LYS:HB2	1.77	0.67
1:D:20:ALA:HA	1:D:120:SER:HA	1.75	0.67
1:D:205:VAL:HA	1:D:208:LEU:HD12	1.75	0.67
1:B:272:ARG:NH1	2:B:395:HOH:O	2.27	0.66
1:B:219:ASN:HD22	1:B:250:TYR:H	1.43	0.66
1:C:61:MET:HE3	1:C:87:HIS:NE2	2.11	0.65
1:B:174:VAL:HB	1:B:275:THR:HG22	1.78	0.65
1:D:205:VAL:HG12	1:D:208:LEU:HD12	1.77	0.65
1:C:77:CYS:SG	1:C:107:LYS:HD2	2.37	0.65
1:A:214:VAL:CG1	1:A:303:ARG:HG3	2.27	0.64
1:B:222:ASN:OD1	1:D:62:ASN:HB2	1.97	0.64
1:B:205:VAL:HG12	1:B:208:LEU:HD12	1.78	0.64
1:B:163:THR:HG22	1:B:170:THR:O	1.98	0.64
1:B:32:VAL:O	1:B:36:VAL:HG23	1.98	0.64
1:A:224:SER:OG	1:A:226:LEU:HB2	1.97	0.63
1:B:69:LEU:O	1:B:73:ILE:HD12	1.98	0.63
1:B:163:THR:HG21	1:B:172:LYS:HD2	1.81	0.62
1:C:313:THR:HG21	1:C:318:GLN:HB3	1.82	0.62
1:D:30:VAL:CG1	1:D:34:ASN:HB2	2.30	0.62
1:C:12:AEI:HZ	1:C:27:VAL:CG2	2.28	0.61
1:B:321:GLN:HG2	1:B:325:GLN:HE21	1.66	0.61
1:A:224:SER:HB2	1:C:236:TYR:OH	2.00	0.61
1:A:12:AEI:HZ	1:A:27:VAL:HG11	1.82	0.61
1:B:269:ARG:O	1:B:278:THR:HG21	2.01	0.61
1:A:69:LEU:O	1:A:73:ILE:HD12	1.99	0.61
1:B:285:ASP:OD2	1:B:288:LYS:HB2	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:72:LYS:HD2	2:C:511:HOH:O	2.01	0.61
1:C:226:LEU:HB2	1:C:227:PRO:HD3	1.83	0.61
1:D:30:VAL:HG13	1:D:34:ASN:HB2	1.82	0.60
1:D:182:ILE:HG12	1:D:187:ILE:HG12	1.83	0.60
1:B:191:ARG:HG3	1:C:191:ARG:HD2	1.82	0.60
1:C:11:GLY:C	1:C:27:VAL:HG13	2.27	0.59
1:A:247:GLY:HA3	1:A:278:THR:HG23	1.83	0.59
1:C:26:THR:OG1	1:C:29:LYS:HE3	2.03	0.59
1:A:60:ASP:OD1	1:C:251:LYS:HB3	2.03	0.59
1:B:40:PRO:HA	2:B:349:HOH:O	2.03	0.58
1:B:142:ALA:O	1:B:143:ASN:HB2	2.02	0.58
1:C:63:ASP:HB3	1:C:211:LEU:HD22	1.85	0.58
1:D:110:VAL:HG13	1:D:149:VAL:HG23	1.84	0.58
1:B:22:LYS:HE3	2:B:342:HOH:O	2.03	0.57
1:A:276:GLY:HA2	2:C:418:HOH:O	2.04	0.57
1:A:243:GLY:O	1:A:272:ARG:HG2	2.03	0.57
1:D:6:ILE:HD12	1:D:48:VAL:HG11	1.86	0.57
1:C:32:VAL:O	1:C:36:VAL:HG23	2.04	0.57
1:B:169:ALA:O	1:B:172:LYS:HE3	2.04	0.57
1:C:244:VAL:HG12	1:C:272:ARG:CZ	2.35	0.57
1:D:67:LEU:HD21	1:D:211:LEU:CD1	2.35	0.56
1:B:97:TYR:CE1	1:B:158:ARG:HG3	2.40	0.56
1:B:205:VAL:HA	1:B:208:LEU:HD12	1.88	0.56
1:D:97:TYR:CE1	1:D:158:ARG:HG3	2.40	0.56
1:A:251:LYS:HD3	1:A:252:SER:N	2.21	0.56
1:A:314:LYS:HE2	2:A:435:HOH:O	2.05	0.55
1:B:65:VAL:O	1:B:69:LEU:HB2	2.06	0.55
1:D:249:LEU:HD23	1:D:257:LEU:HD12	1.88	0.55
1:C:205:VAL:HG12	1:C:208:LEU:HD12	1.89	0.55
1:C:65:VAL:HG12	1:C:69:LEU:HD23	1.88	0.55
1:D:225:ASP:HB3	1:D:252:SER:HB2	1.88	0.55
1:A:27:VAL:HG21	1:A:57:GLY:HA3	1.88	0.54
1:A:96:ALA:O	1:A:100:ASP:HB2	2.07	0.54
1:D:67:LEU:HD21	1:D:211:LEU:HD11	1.90	0.54
1:A:225:ASP:O	1:A:229:LYS:HG3	2.08	0.54
1:A:7:LEU:HD22	1:A:72:LYS:HD3	1.90	0.54
1:C:3:ASN:O	1:C:81:ASP:HB2	2.08	0.53
1:A:11:GLY:HA2	1:A:12:AEI:OT1	2.09	0.52
1:B:219:ASN:ND2	1:B:250:TYR:H	2.06	0.52
1:B:321:GLN:OE1	1:B:325:GLN:NE2	2.42	0.52
1:B:12:AEI:HZ	1:B:27:VAL:HG11	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:15:GLY:O	1:D:118:SER:OG	2.27	0.52
1:B:159:ASP:O	1:B:173:SER:HA	2.09	0.52
2:B:456:HOH:O	1:C:188:ASP:HB3	2.10	0.52
1:A:205:VAL:HG22	2:A:346:HOH:O	2.08	0.52
1:D:251:LYS:HE2	1:D:255:ASP:OD2	2.10	0.52
1:A:169:ALA:O	1:A:172:LYS:HE3	2.09	0.52
1:B:27:VAL:HG21	1:B:57:GLY:CA	2.40	0.51
1:D:62:ASN:N	1:D:62:ASN:HD22	2.09	0.51
1:A:12:AEI:HB	1:A:114:ALA:O	2.11	0.51
1:C:17:GLY:HA3	1:C:24:ASN:O	2.09	0.51
1:C:303:ARG:HD3	2:C:335:HOH:O	2.11	0.51
1:B:19:SER:HB3	1:B:22:LYS:HB3	1.93	0.51
1:C:69:LEU:O	1:C:73:ILE:HG13	2.09	0.51
1:A:286:ASP:HB3	1:A:291:PHE:O	2.11	0.51
1:B:163:THR:HG21	1:B:172:LYS:NZ	2.26	0.51
1:C:110:VAL:HG13	1:C:149:VAL:HG23	1.92	0.51
1:D:12:AEI:HZ	1:D:27:VAL:HG11	1.93	0.51
1:C:75:THR:HG22	1:C:76:ASP:OD1	2.11	0.51
1:D:219:ASN:HD22	1:D:250:TYR:H	1.57	0.51
1:C:35:LEU:O	1:C:39:VAL:HG23	2.09	0.51
1:D:66:TRP:HB3	1:D:98:PHE:CE2	2.46	0.51
1:A:8:ALA:HA	1:A:86:THR:OG1	2.11	0.50
1:A:292:VAL:HG22	1:A:320:GLN:HB2	1.91	0.50
1:B:19:SER:HB3	1:B:22:LYS:CB	2.41	0.50
1:D:320:GLN:O	1:D:324:ASN:OD1	2.29	0.50
1:B:11:GLY:HA2	1:B:12:AEI:OT2	2.11	0.50
1:D:24:ASN:O	1:D:25:TYR:HB3	2.12	0.50
1:D:115:MET:HE2	1:D:170:THR:HG21	1.94	0.50
1:C:220:TYR:CE2	1:C:223:ALA:HA	2.47	0.50
1:C:21:THR:HG22	1:C:121:MET:CE	2.42	0.50
1:A:190:GLN:HA	1:D:194:ALA:HB3	1.94	0.50
1:A:236:TYR:OH	1:C:224:SER:HB2	2.11	0.49
1:A:214:VAL:HG12	1:A:303:ARG:HG3	1.94	0.49
1:A:216:ILE:HA	1:A:240:VAL:O	2.12	0.49
1:B:61:MET:HE3	1:B:87:HIS:NE2	2.27	0.49
2:A:355:HOH:O	1:C:272:ARG:HD3	2.10	0.49
1:D:6:ILE:HD12	1:D:48:VAL:CG1	2.43	0.49
1:A:33:GLU:CD	1:A:33:GLU:H	2.18	0.49
1:D:174:VAL:HB	1:D:275:THR:HG22	1.95	0.49
1:C:182:ILE:HG12	1:C:187:ILE:HG12	1.93	0.49
1:D:58:SER:HB3	1:D:87:HIS:CE1	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:ASN:HB3	1:C:303:ARG:NH2	2.28	0.49
1:C:67:LEU:HD21	1:C:211:LEU:CD1	2.43	0.49
1:D:228:ALA:O	1:D:232:VAL:HG23	2.12	0.49
1:A:67:LEU:HD21	1:A:211:LEU:HD11	1.95	0.48
1:A:220:TYR:HA	1:A:272:ARG:NH2	2.28	0.48
1:C:8:ALA:HA	1:C:86:THR:OG1	2.13	0.48
1:D:314:LYS:O	1:D:316:PRO:HD3	2.13	0.48
1:D:147:LEU:HG	1:D:154:VAL:CG1	2.44	0.48
1:A:285:ASP:O	1:A:289:TYR:CD2	2.67	0.48
1:A:14:ALA:HA	1:A:30:VAL:HB	1.96	0.48
1:C:109:VAL:HG12	1:C:146:VAL:HG13	1.96	0.47
1:A:39:VAL:HB	1:A:42:LEU:HD22	1.95	0.47
1:A:58:SER:HB2	1:A:91:THR:OG1	2.13	0.47
1:D:32:VAL:O	1:D:36:VAL:HG23	2.14	0.47
1:D:247:GLY:HA3	1:D:278:THR:HG23	1.97	0.47
1:C:214:VAL:CG1	1:C:303:ARG:HG3	2.44	0.47
1:A:84:VAL:HA	1:A:110:VAL:O	2.14	0.47
1:C:116:ARG:HB3	1:C:120:SER:OG	2.14	0.47
1:D:196:LYS:HB2	1:D:326:TYR:CD1	2.49	0.47
1:B:42:LEU:HD11	1:B:130:TYR:HA	1.96	0.47
1:C:249:LEU:HD23	1:C:254:PHE:HA	1.95	0.47
1:B:204:ASP:HA	2:B:355:HOH:O	2.15	0.47
1:C:66:TRP:HB3	1:C:98:PHE:CE2	2.50	0.47
1:A:163:THR:O	1:C:274:PRO:HG2	2.14	0.46
1:C:49:LYS:HE2	1:C:51:GLU:OE2	2.15	0.46
1:A:248:ASN:HA	2:A:390:HOH:O	2.15	0.46
1:C:99:LEU:O	1:C:103:VAL:HB	2.14	0.46
1:D:249:LEU:CD2	1:D:257:LEU:HD12	2.45	0.46
1:C:61:MET:HE3	1:C:87:HIS:CD2	2.50	0.46
1:B:214:VAL:CG1	1:B:306:LEU:HD23	2.45	0.46
1:C:241:SER:HB3	2:C:395:HOH:O	2.15	0.46
1:D:178:PRO:HG3	1:D:181:TYR:OH	2.16	0.46
1:A:155:LEU:HD13	1:A:160:VAL:HG23	1.98	0.46
1:D:100:ASP:HA	1:D:146:VAL:HG23	1.97	0.46
1:A:222:ASN:HB2	2:C:334:HOH:O	2.15	0.46
1:B:273:VAL:HG11	1:D:164:ASN:C	2.40	0.46
1:C:219:ASN:ND2	1:C:250:TYR:H	2.13	0.46
1:A:213:LYS:HE3	1:A:236:TYR:CE1	2.51	0.46
1:A:90:ASP:HB3	1:C:244:VAL:HB	1.97	0.46
1:A:100:ASP:HA	1:A:146:VAL:HG23	1.97	0.46
1:A:216:ILE:HG12	1:A:240:VAL:HB	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:224:SER:OG	1:C:226:LEU:HG	2.16	0.45
1:A:28:GLY:HA2	1:A:55:ASN:OD1	2.17	0.45
1:B:162:LYS:NZ	1:D:245:GLY:HA2	2.31	0.45
1:C:247:GLY:HA3	1:C:278:THR:HG23	1.99	0.45
1:D:67:LEU:CD2	1:D:205:VAL:HB	2.45	0.45
1:A:89:VAL:HG21	1:A:115:MET:SD	2.57	0.45
1:A:251:LYS:HD3	1:A:252:SER:H	1.82	0.45
1:B:272:ARG:NH1	2:D:344:HOH:O	2.49	0.45
1:C:61:MET:SD	1:C:65:VAL:HG11	2.58	0.45
1:A:42:LEU:HA	1:A:42:LEU:HD12	1.74	0.44
1:B:8:ALA:HB3	1:B:52:GLN:HA	1.98	0.44
1:B:226:LEU:HD21	1:D:234:ALA:HB2	1.98	0.44
1:D:214:VAL:HA	1:D:238:GLY:O	2.16	0.44
1:D:286:ASP:O	1:D:290:GLY:N	2.48	0.44
1:B:99:LEU:O	1:B:103:VAL:HB	2.16	0.44
1:B:174:VAL:HB	1:B:275:THR:CG2	2.47	0.44
1:A:191:ARG:NH2	1:D:156:ASP:OD1	2.51	0.44
1:A:270:SER:HA	1:A:294:SER:OG	2.17	0.44
1:B:182:ILE:HG12	1:B:187:ILE:HG12	1.99	0.44
1:C:65:VAL:O	1:C:69:LEU:HD23	2.17	0.44
1:C:212:PRO:O	1:C:307:GLN:HG2	2.17	0.44
1:A:129:LEU:O	1:A:130:TYR:C	2.57	0.44
1:A:285:ASP:O	1:A:289:TYR:HD2	1.99	0.44
1:C:288:LYS:O	1:C:288:LYS:HG3	2.17	0.44
1:D:174:VAL:CG2	1:D:275:THR:HG22	2.47	0.44
1:A:14:ALA:HA	1:A:35:LEU:HD21	2.00	0.44
1:A:212:PRO:O	1:A:307:GLN:HG2	2.16	0.44
1:B:315:ASP:HB3	1:B:318:GLN:HB2	2.00	0.44
1:C:249:LEU:HD12	1:C:249:LEU:N	2.32	0.44
1:A:16:GLY:HA2	1:A:119:THR:HG23	2.00	0.44
1:B:156:ASP:OD2	1:C:191:ARG:NH2	2.51	0.44
1:B:174:VAL:CB	1:B:275:THR:HG22	2.47	0.44
1:B:213:LYS:HD3	2:B:387:HOH:O	2.18	0.44
1:B:226:LEU:CB	1:B:227:PRO:HD3	2.48	0.44
1:D:292:VAL:HG13	1:D:320:GLN:HA	2.00	0.44
1:B:321:GLN:HG2	1:B:325:GLN:NE2	2.33	0.43
1:C:115:MET:HE2	1:C:170:THR:HG21	2.00	0.43
1:C:89:VAL:HG11	1:C:115:MET:SD	2.58	0.43
1:D:268:VAL:CG1	1:D:294:SER:HB3	2.48	0.43
1:A:78:ASP:OD1	1:A:79:LYS:HG3	2.18	0.43
1:B:20:ALA:HA	1:B:120:SER:HA	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:178:PRO:HG3	1:D:181:TYR:CZ	2.53	0.43
1:D:206:SER:OG	1:D:207:LYS:HD2	2.18	0.43
1:A:10:GLY:HA3	1:A:56:ILE:O	2.19	0.43
1:B:191:ARG:CG	1:C:191:ARG:HD2	2.48	0.43
1:C:220:TYR:CZ	1:C:223:ALA:HA	2.53	0.43
1:D:121:MET:HB2	1:D:121:MET:HE3	1.72	0.43
1:A:1:LEU:N	2:A:447:HOH:O	2.42	0.43
1:B:220:TYR:CZ	1:B:223:ALA:HA	2.54	0.43
1:D:69:LEU:O	1:D:73:ILE:HG13	2.18	0.43
1:D:76:ASP:O	1:D:77:CYS:C	2.62	0.43
1:B:272:ARG:NH2	2:B:462:HOH:O	2.42	0.43
1:B:301:LYS:O	1:B:304:VAL:HB	2.18	0.43
1:C:12:AEI:HB	1:C:114:ALA:O	2.18	0.43
1:D:315:ASP:OD2	1:D:318:GLN:HG3	2.19	0.43
1:C:286:ASP:O	1:C:290:GLY:N	2.50	0.43
1:D:142:ALA:O	1:D:143:ASN:HB2	2.18	0.43
1:D:269:ARG:O	1:D:278:THR:HG21	2.19	0.43
1:A:101:LEU:CD1	1:A:308:LEU:HD12	2.49	0.43
1:A:171:PHE:C	1:A:172:LYS:HG3	2.44	0.43
1:B:214:VAL:HG13	1:B:306:LEU:HD23	2.00	0.43
1:C:286:ASP:HB3	1:C:291:PHE:O	2.18	0.43
1:A:65:VAL:HG12	1:A:69:LEU:HD22	2.01	0.42
1:A:66:TRP:HB3	1:A:98:PHE:CE2	2.54	0.42
1:C:11:GLY:CA	1:C:27:VAL:HG13	2.49	0.42
1:D:31:GLY:O	1:D:35:LEU:HB2	2.19	0.42
1:D:96:ALA:O	1:D:100:ASP:HB2	2.19	0.42
1:B:156:ASP:HB3	1:B:179:LEU:HD21	2.00	0.42
1:B:163:THR:HG21	1:B:172:LYS:CD	2.49	0.42
1:C:219:ASN:HD22	1:C:250:TYR:H	1.67	0.42
1:C:254:PHE:CD1	1:C:254:PHE:C	2.97	0.42
1:B:320:GLN:HE21	1:B:324:ASN:HD21	1.68	0.42
1:A:27:VAL:HG21	1:A:57:GLY:CA	2.49	0.42
1:D:8:ALA:HA	1:D:86:THR:OG1	2.19	0.42
1:A:67:LEU:HD23	1:A:67:LEU:HA	1.92	0.42
1:A:214:VAL:HG13	1:A:306:LEU:HD23	2.02	0.42
1:A:244:VAL:HG12	1:A:272:ARG:CZ	2.50	0.42
1:C:96:ALA:O	1:C:100:ASP:HB2	2.19	0.42
1:D:151:ASN:O	1:D:152:ASP:HB2	2.19	0.42
1:A:234:ALA:HB2	1:C:226:LEU:HD13	2.02	0.42
1:D:39:VAL:CG1	1:D:42:LEU:HD13	2.50	0.42
1:B:58:SER:HB2	1:B:91:THR:OG1	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:354:HOH:O	1:D:222:ASN:HB2	2.20	0.42
1:A:226:LEU:HD12	1:A:226:LEU:HA	1.82	0.42
1:B:69:LEU:HB3	1:B:99:LEU:HD21	2.01	0.42
1:C:213:LYS:HE2	2:C:439:HOH:O	2.19	0.42
1:A:231:LEU:O	1:A:236:TYR:HB2	2.20	0.41
1:B:257:LEU:HB3	1:B:267:VAL:HG11	2.01	0.41
1:C:238:GLY:HA2	1:C:266:ALA:O	2.20	0.41
1:A:105:CYS:SG	1:A:107:LYS:HB2	2.60	0.41
1:A:130:TYR:O	1:A:134:VAL:HG23	2.20	0.41
1:B:261:ALA:HA	1:B:265:THR:O	2.20	0.41
1:C:218:TYR:HA	1:C:242:ALA:HB3	2.03	0.41
1:A:6:ILE:HA	1:A:84:VAL:O	2.19	0.41
1:C:106:ASP:HB3	1:C:143:ASN:OD1	2.21	0.41
1:D:159:ASP:OD1	1:D:175:ASN:HB2	2.20	0.41
1:D:286:ASP:HB3	1:D:291:PHE:O	2.20	0.41
1:C:67:LEU:HD21	1:C:211:LEU:HD11	2.02	0.41
1:C:94:GLU:CD	1:C:300:GLN:HB2	2.46	0.41
1:C:214:VAL:HG12	1:C:303:ARG:HG3	2.01	0.41
1:D:139:LYS:HD3	1:D:139:LYS:HA	1.84	0.41
1:D:196:LYS:HD3	1:D:326:TYR:CE1	2.56	0.41
1:D:243:GLY:O	1:D:272:ARG:HG2	2.19	0.41
1:B:286:ASP:HB2	2:B:408:HOH:O	2.20	0.41
1:B:163:THR:HG22	1:B:170:THR:HA	2.02	0.41
1:B:241:SER:HB3	2:B:393:HOH:O	2.20	0.41
1:A:219:ASN:ND2	1:A:250:TYR:H	2.18	0.41
1:A:286:ASP:O	1:A:290:GLY:N	2.50	0.41
1:D:72:LYS:HB2	1:D:72:LYS:HE2	1.76	0.41
1:A:315:ASP:HA	1:A:316:PRO:HD2	1.95	0.41
1:B:66:TRP:HB3	1:B:98:PHE:CD2	2.56	0.41
1:B:84:VAL:HA	1:B:110:VAL:O	2.20	0.41
1:B:131:ASN:HD22	1:B:131:ASN:HA	1.75	0.41
1:B:226:LEU:HB2	1:B:227:PRO:HD3	2.03	0.41
1:B:260:ALA:O	1:B:265:THR:HB	2.20	0.41
1:B:319:ILE:O	1:B:323:PHE:HD1	2.04	0.41
1:C:67:LEU:HD21	1:C:211:LEU:HD13	2.03	0.41
1:D:194:ALA:HB2	2:D:338:HOH:O	2.19	0.41
1:D:249:LEU:HD22	1:D:254:PHE:HA	2.02	0.41
1:D:249:LEU:N	1:D:249:LEU:HD12	2.36	0.41
1:A:98:PHE:CE1	1:A:211:LEU:HD21	2.55	0.41
1:C:266:ALA:HB2	1:C:310:LEU:HD21	2.02	0.41
1:A:107:LYS:O	1:A:109:VAL:HG23	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:220:TYR:HA	1:D:272:ARG:NH2	2.36	0.40
1:B:190:GLN:O	1:B:191:ARG:CD	2.69	0.40
1:C:97:TYR:CE1	1:C:158:ARG:HG3	2.57	0.40
1:D:99:LEU:O	1:D:103:VAL:HB	2.21	0.40
1:C:248:ASN:C	1:C:249:LEU:HD12	2.46	0.40
1:B:312:GLN:NE2	2:B:442:HOH:O	2.55	0.40
1:A:171:PHE:O	1:A:172:LYS:HG3	2.22	0.40
1:A:247:GLY:HA2	1:A:270:SER:O	2.21	0.40
1:D:7:LEU:O	1:D:85:ILE:HA	2.21	0.40
1:D:272:ARG:HG3	1:D:273:VAL:N	2.37	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	323/326 (99%)	311 (96%)	10 (3%)	2 (1%)	21 23
1	B	323/326 (99%)	305 (94%)	18 (6%)	0	100 100
1	C	323/326 (99%)	306 (95%)	16 (5%)	1 (0%)	36 42
1	D	323/326 (99%)	307 (95%)	16 (5%)	0	100 100
All	All	1292/1304 (99%)	1229 (95%)	60 (5%)	3 (0%)	43 51

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	251	LYS
1	C	126	PRO
1	A	126	PRO

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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	AEI	C	12	1	12,14,15	1.15	2 (16%)	12,18,20	1.49	2 (16%)
1	AEI	B	12	1	12,14,15	1.19	1 (8%)	12,18,20	1.50	2 (16%)
1	AEI	A	12	1	12,14,15	1.39	2 (16%)	12,18,20	1.73	3 (25%)
1	AEI	D	12	1	12,14,15	1.15	1 (8%)	12,18,20	1.39	2 (16%)

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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	AEI	C	12	1	-	4/16/18/20	-
1	AEI	B	12	1	-	5/16/18/20	-
1	AEI	A	12	1	-	4/16/18/20	-
1	AEI	D	12	1	-	6/16/18/20	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	12	AEI	CB-CA	2.70	1.59	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	12	AEI	OG1-CD	2.48	1.41	1.34
1	C	12	AEI	OG1-CD	2.18	1.40	1.34
1	C	12	AEI	OT2-CH2	-2.12	1.23	1.30
1	D	12	AEI	OG1-CB	-2.10	1.43	1.46
1	B	12	AEI	OT2-CH2	-2.09	1.24	1.30

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	12	AEI	OG1-CD-CE2	4.22	118.99	111.43
1	B	12	AEI	OG1-CD-CE2	3.60	117.88	111.43
1	D	12	AEI	OG1-CD-CE2	3.58	117.85	111.43
1	C	12	AEI	OG1-CD-CE2	3.52	117.75	111.43
1	A	12	AEI	OE1-CD-CE2	-2.95	117.80	124.65
1	B	12	AEI	OE1-CD-CE2	-2.44	118.97	124.65
1	C	12	AEI	OE1-CD-CE2	-2.43	119.00	124.65
1	A	12	AEI	OG1-CB-CA	2.15	110.53	105.80
1	D	12	AEI	OE1-CD-CE2	-2.07	119.83	124.65

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	12	AEI	O-C-CA-CB
1	B	12	AEI	O-C-CA-CB
1	C	12	AEI	O-C-CA-CB
1	D	12	AEI	O-C-CA-CB
1	D	12	AEI	OT2-CH2-CZ-NH1
1	B	12	AEI	OT2-CH2-CZ-NH1
1	C	12	AEI	OT2-CH2-CZ-NH1
1	B	12	AEI	OT1-CH2-CZ-NH1
1	D	12	AEI	OT1-CH2-CZ-NH1
1	A	12	AEI	OE1-CD-CE2-CZ
1	B	12	AEI	OE1-CD-CE2-CZ
1	A	12	AEI	OG1-CD-CE2-CZ
1	B	12	AEI	OG1-CD-CE2-CZ
1	A	12	AEI	OT2-CH2-CZ-NH1
1	D	12	AEI	OG1-CD-CE2-CZ
1	C	12	AEI	OE1-CD-CE2-CZ
1	D	12	AEI	CD-CE2-CZ-CH2
1	D	12	AEI	OE1-CD-CE2-CZ
1	C	12	AEI	OT1-CH2-CZ-NH1

There are no ring outliers.

4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	12	AEI	3	0
1	B	12	AEI	2	0
1	A	12	AEI	3	0
1	D	12	AEI	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	325/326 (99%)	-0.75	1 (0%) 90 88	5, 17, 39, 58	0
1	B	325/326 (99%)	-0.72	1 (0%) 90 88	5, 17, 42, 66	0
1	C	325/326 (99%)	-0.77	1 (0%) 90 88	5, 17, 42, 62	0
1	D	325/326 (99%)	-0.76	0 100 100	6, 16, 41, 59	0
All	All	1300/1304 (99%)	-0.75	3 (0%) 91 90	5, 17, 41, 66	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	321	GLN	2.7
1	C	18	ASP	2.5
1	B	321	GLN	2.4

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	AEI	D	12	15/16	0.94	0.07	14,22,28,30	0
1	AEI	C	12	15/16	0.96	0.06	16,26,28,30	0
1	AEI	A	12	15/16	0.96	0.06	11,21,27,28	0
1	AEI	B	12	15/16	0.97	0.06	17,23,26,27	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.