



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:53 AM UTC

PDB ID : 1K28 / pdb_00001k28
Title : The Structure of the Bacteriophage T4 Cell-Puncturing Device
Authors : Kanamaru, S.; Leiman, P.G.; Kostyuchenko, V.A.; Chipman, P.R.; Mesyanzhinov, V.V.; Arisaka, F.; Rossmann, M.G.
Deposited on : 2001-09-26
Resolution : 2.90 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

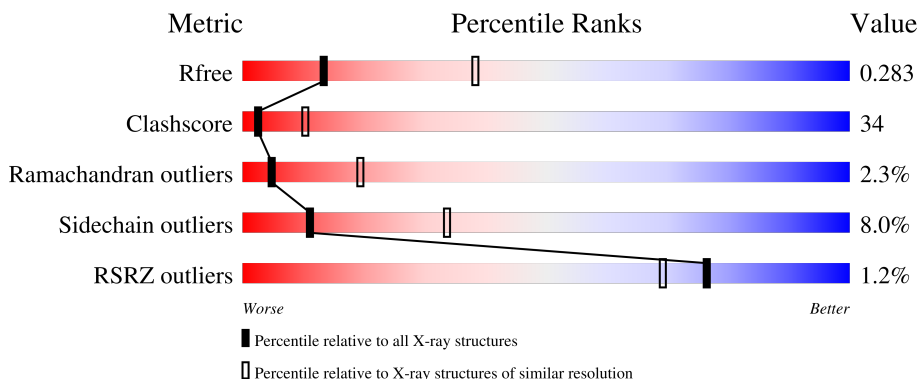
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	584	
2	D	391	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7614 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 TAIL-ASSOCIATED LYSOZYME.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	557	Total	C	N	O	S	0	0	0
			4309	2672	763	853	21			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	576	SER	-	expression tag	UNP P16009
A	577	VAL	-	expression tag	UNP P16009
A	578	ASP	-	expression tag	UNP P16009
A	579	HIS	-	expression tag	UNP P16009
A	580	HIS	-	expression tag	UNP P16009
A	581	HIS	-	expression tag	UNP P16009
A	582	HIS	-	expression tag	UNP P16009
A	583	HIS	-	expression tag	UNP P16009
A	584	HIS	-	expression tag	UNP P16009

- Molecule 2 is a protein called BASEPLATE STRUCTURAL PROTEIN GP27.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	D	364	Total	C	N	O	S	Se	0	0	0
			2912	1858	478	559	2	15			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	1	MSE	MET	modified residue	UNP P17172
D	3	MSE	MET	modified residue	UNP P17172
D	40	MSE	MET	modified residue	UNP P17172
D	51	MSE	MET	modified residue	UNP P17172
D	64	MSE	MET	modified residue	UNP P17172
D	135	MSE	MET	modified residue	UNP P17172
D	168	ASN	THR	SEE REMARK 999	UNP P17172

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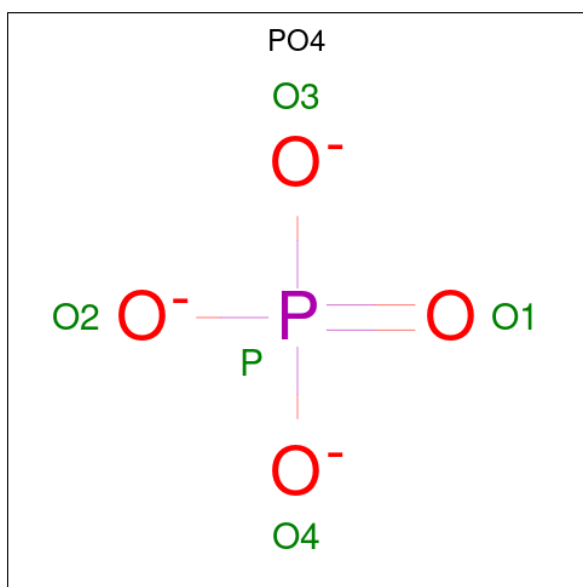
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Chain	Residue	Modelled	Actual	Comment	Reference
D	193	MSE	MET	modified residue	UNP P17172
D	197	MSE	MET	modified residue	UNP P17172
D	198	MSE	MET	modified residue	UNP P17172
D	202	MSE	MET	modified residue	UNP P17172
D	203	MSE	MET	modified residue	UNP P17172
D	211	MSE	MET	modified residue	UNP P17172
D	248	MSE	MET	modified residue	UNP P17172
D	262	SER	GLU	SEE REMARK 999	UNP P17172
D	265	MSE	MET	modified residue	UNP P17172
D	286	MSE	MET	modified residue	UNP P17172
D	301	MSE	MET	modified residue	UNP P17172
D	308	ALA	LEU	SEE REMARK 999	UNP P17172
D	357	MSE	MET	modified residue	UNP P17172

- Molecule 3 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total K 1 1	0	0

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O P 3 2 1	0	0

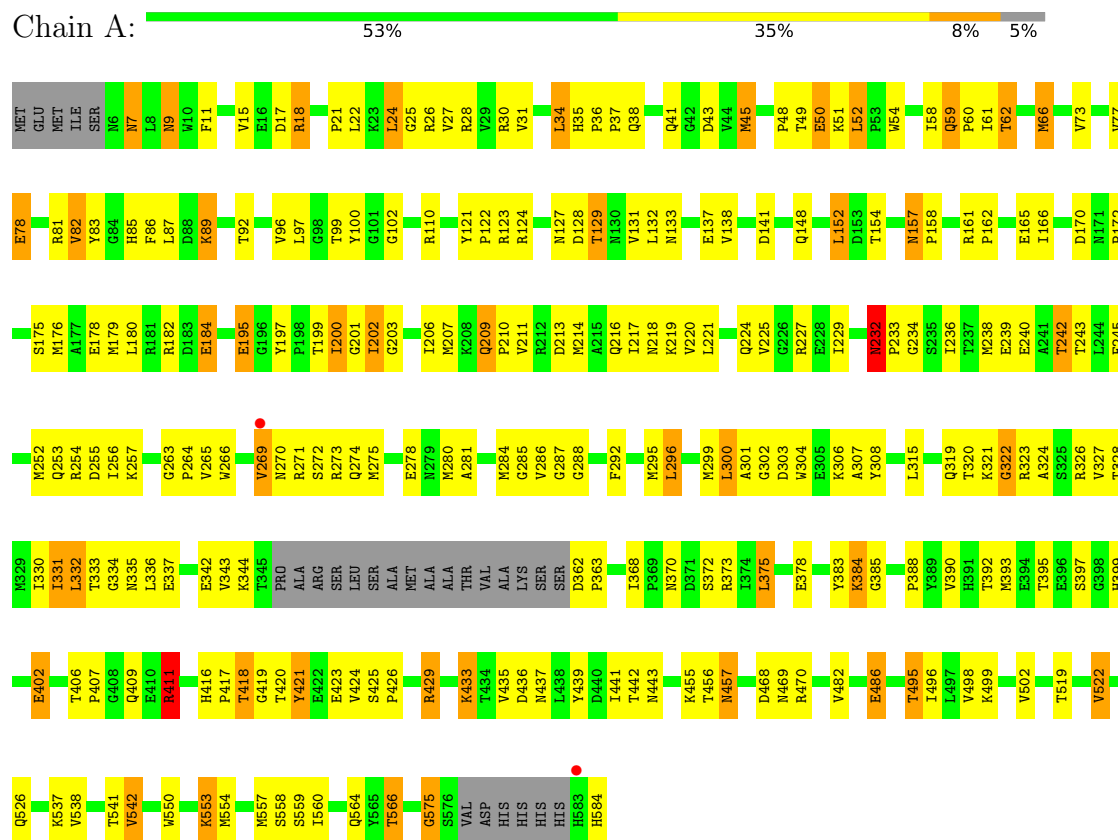
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	307	Total 307	O 307	0	0
5	D	82	Total 82	O 82	0	0

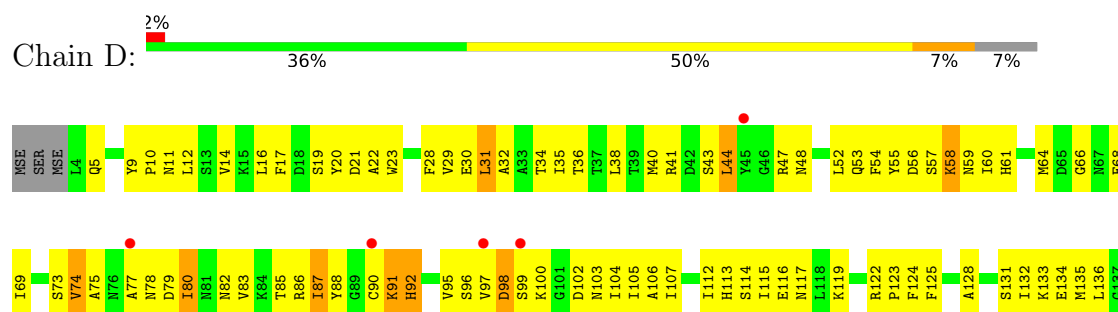
3 Residue-property plots

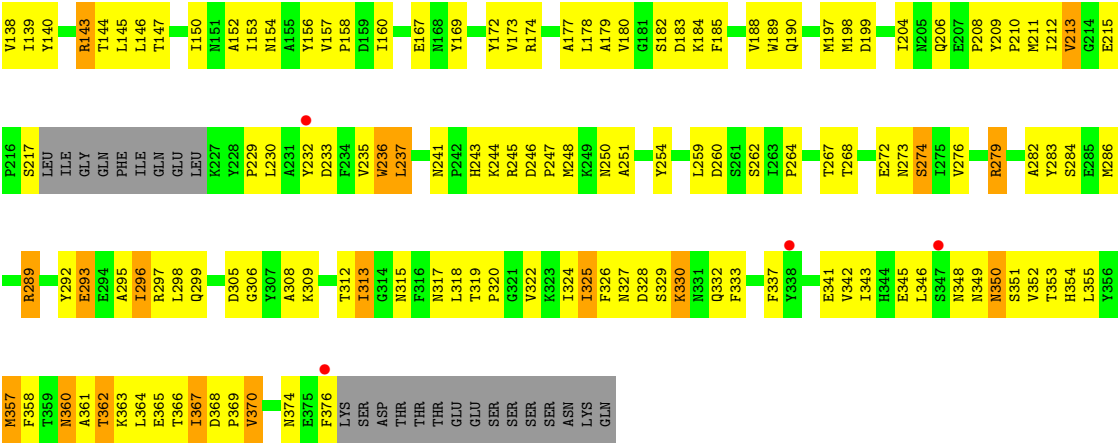
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: TAIL-ASSOCIATED LYSOZYME



• Molecule 2: BASEPLATE STRUCTURAL PROTEIN GP27





4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	139.26Å 139.26Å 382.00Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.90 50.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	5.1 (50.00-2.90) 98.9 (50.00-2.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.57 (at 2.91Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.211 , 0.280 0.215 , 0.283	Depositor DCC
R_{free} test set	1602 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	52.2	Xtriage
Anisotropy	0.794	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 79.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7614	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.73% 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: K, PO4

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	0.47	0/4391	0.99	18/5950 (0.3%)
2	D	0.38	0/2965	0.90	6/3997 (0.2%)
All	All	0.44	0/7356	0.95	24/9947 (0.2%)

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	232	ASN	CA-C-N	-8.00	109.85	119.84
1	A	232	ASN	C-N-CA	-8.00	109.85	119.84
1	A	157	ASN	CA-C-N	6.46	126.49	119.90
1	A	157	ASN	C-N-CA	6.46	126.49	119.90
1	A	443	ASN	N-CA-C	6.22	118.14	111.36
1	A	421	TYR	N-CA-C	6.04	117.37	108.86
1	A	269	VAL	N-CA-C	5.86	118.88	110.09
1	A	402	GLU	N-CA-C	5.77	118.31	108.90
1	A	411	ARG	N-CA-C	5.72	117.55	108.79
2	D	78	ASN	N-CA-C	5.52	116.97	111.07
1	A	575	GLY	N-CA-C	5.50	126.21	113.18
1	A	82	VAL	N-CA-C	5.49	118.01	108.95
2	D	44	LEU	N-CA-C	-5.49	106.54	113.18
1	A	158	PRO	N-CA-C	5.33	119.56	111.15
1	A	161	ARG	CA-C-N	-5.32	114.48	119.85
1	A	161	ARG	C-N-CA	-5.32	114.48	119.85
1	A	166	ILE	CA-C-N	5.27	125.19	119.76
1	A	166	ILE	C-N-CA	5.27	125.19	119.76
2	D	143	ARG	N-CA-C	-5.22	104.71	111.71
1	A	560	ILE	N-CA-C	5.19	115.97	107.24
2	D	279	ARG	N-CA-C	-5.19	104.24	110.88
2	D	274	SER	N-CA-C	5.18	117.27	109.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	86	PHE	N-CA-C	-5.03	100.97	109.07
2	D	284	SER	N-CA-C	-5.02	107.10	113.18

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	4309	0	4202	257	1
2	D	2912	0	2835	237	0
3	A	1	0	0	0	0
4	A	3	0	0	0	1
5	A	307	0	0	6	0
5	D	82	0	0	5	0
All	All	7614	0	7037	489	2

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 34.

All (489) 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:89:LYS:HE2	1:A:89:LYS:H	1.16	1.10
1:A:184:GLU:HG2	1:A:203:GLY:HA3	1.36	1.03
1:A:89:LYS:H	1:A:89:LYS:CE	1.72	1.03
1:A:221:LEU:HD11	1:A:236:ILE:HD11	1.40	1.01
1:A:416:HIS:CD2	1:A:418:THR:HB	1.95	1.01
1:A:330:ILE:HD12	1:A:336:LEU:HD12	1.43	0.99
1:A:416:HIS:HD2	1:A:418:THR:HB	1.27	0.98
1:A:89:LYS:HE2	1:A:89:LYS:N	1.78	0.97
2:D:330:LYS:HA	2:D:330:LYS:HE3	1.52	0.91
2:D:95:VAL:HA	2:D:105:ILE:HG22	1.54	0.90
1:A:127:ASN:ND2	1:A:129:THR:HG23	1.87	0.90
1:A:127:ASN:HD21	1:A:129:THR:HG23	1.37	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:315:ASN:HD21	2:D:317:ASN:HD22	1.19	0.88
1:A:307:ALA:HB1	1:A:331:ILE:HD11	1.55	0.88
2:D:286:MSE:HE3	2:D:286:MSE:HA	1.56	0.88
1:A:221:LEU:HD12	1:A:229:ILE:HG13	1.53	0.87
1:A:256:ILE:HA	1:A:286:VAL:HG21	1.55	0.87
1:A:31:VAL:HB	1:A:34:LEU:HD11	1.57	0.86
1:A:133:ASN:HD21	1:A:392:THR:H	1.22	0.84
2:D:11:ASN:H	2:D:74:VAL:HG23	1.41	0.83
1:A:48:PRO:HG2	1:A:51:LYS:HD3	1.62	0.81
1:A:270:ASN:ND2	1:A:271:ARG:H	1.76	0.81
1:A:31:VAL:HB	1:A:34:LEU:CD1	2.11	0.81
2:D:75:ALA:HB2	2:D:83:VAL:HG12	1.65	0.79
2:D:87:ILE:HD12	2:D:87:ILE:H	1.45	0.79
2:D:79:ASP:OD1	2:D:82:ASN:HB3	1.83	0.79
2:D:11:ASN:N	2:D:74:VAL:HG23	1.97	0.79
1:A:266:TRP:NE1	1:A:274:GLN:HG2	1.99	0.78
2:D:169:TYR:O	2:D:173:VAL:HG23	1.83	0.78
2:D:188:VAL:HG12	2:D:197:MSE:HG2	1.67	0.76
2:D:54:PHE:O	2:D:104:ILE:HD12	1.86	0.76
2:D:315:ASN:HD21	2:D:317:ASN:ND2	1.83	0.76
2:D:16:LEU:HD22	2:D:31:LEU:HD21	1.66	0.75
2:D:325:ILE:HD13	2:D:326:PHE:N	2.02	0.75
1:A:221:LEU:O	1:A:225:VAL:HG12	1.87	0.74
2:D:237:LEU:HD22	2:D:237:LEU:H	1.51	0.74
1:A:38:GLN:OE1	1:A:45:MET:HE3	1.88	0.73
2:D:36:THR:HB	2:D:53:GLN:HG3	1.69	0.73
1:A:416:HIS:HD2	1:A:418:THR:H	1.37	0.72
1:A:172:PRO:HG3	1:A:272:SER:OG	1.90	0.72
2:D:96:SER:OG	2:D:104:ILE:HG23	1.90	0.72
2:D:212:ILE:HG22	2:D:327:ASN:HB2	1.72	0.72
2:D:251:ALA:HB2	2:D:273:ASN:HB2	1.72	0.72
1:A:214:MET:HG2	1:A:232:ASN:O	1.90	0.71
1:A:284:MET:HE1	1:A:292:PHE:HE1	1.54	0.71
2:D:80:ILE:HD13	2:D:80:ILE:H	1.55	0.71
2:D:10:PRO:HB2	2:D:74:VAL:HG21	1.73	0.71
2:D:11:ASN:HD21	2:D:80:ILE:HG22	1.55	0.71
1:A:127:ASN:HD22	1:A:128:ASP:N	1.87	0.70
1:A:209:GLN:HG3	1:A:210:PRO:HD2	1.72	0.70
1:A:200:ILE:HG12	1:A:201:GLY:H	1.55	0.70
2:D:134:GLU:O	2:D:138:VAL:HG23	1.92	0.70
2:D:337:PHE:HB2	2:D:357:MSE:HE2	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:ASN:O	1:A:234:GLY:N	2.25	0.69
1:A:416:HIS:CD2	1:A:418:THR:H	2.09	0.69
1:A:176:MET:HE3	1:A:180:LEU:HD21	1.75	0.69
1:A:418:THR:HG22	1:A:420:THR:H	1.56	0.69
2:D:208:PRO:O	2:D:210:PRO:HD3	1.93	0.69
2:D:360:ASN:HD22	2:D:360:ASN:H	1.41	0.69
1:A:470:ARG:HG3	1:A:470:ARG:HH11	1.58	0.68
1:A:409:GLN:HE21	1:A:409:GLN:HA	1.57	0.68
1:A:184:GLU:CG	1:A:203:GLY:HA3	2.18	0.67
1:A:273:ARG:NH1	1:A:334:GLY:HA3	2.09	0.67
2:D:135:MSE:O	2:D:139:ILE:HG12	1.94	0.67
2:D:324:ILE:HD12	2:D:357:MSE:HE3	1.77	0.67
1:A:232:ASN:HB3	1:A:233:PRO:CD	2.25	0.67
1:A:239:GLU:O	1:A:243:THR:HG23	1.95	0.67
1:A:370:ASN:HD21	1:A:372:SER:CB	2.08	0.66
1:A:344:LYS:HA	1:A:344:LYS:HE2	1.78	0.66
1:A:330:ILE:CD1	1:A:336:LEU:HD12	2.24	0.66
2:D:325:ILE:HD13	2:D:326:PHE:H	1.60	0.66
2:D:250:ASN:HD21	2:D:272:GLU:HB2	1.62	0.65
1:A:37:PRO:HG2	1:A:38:GLN:NE2	2.12	0.65
1:A:221:LEU:HD11	1:A:236:ILE:CD1	2.22	0.65
1:A:418:THR:HG23	1:A:435:VAL:HG21	1.78	0.65
2:D:133:LYS:HG2	2:D:150:ILE:HG21	1.79	0.65
1:A:176:MET:O	1:A:180:LEU:HD23	1.97	0.64
2:D:360:ASN:HD22	2:D:360:ASN:N	1.94	0.64
2:D:14:VAL:HG11	2:D:52:LEU:HD11	1.77	0.64
2:D:367:ILE:HG22	2:D:367:ILE:O	1.97	0.64
1:A:495:THR:O	1:A:496:ILE:HD13	1.97	0.64
1:A:308:TYR:HB2	1:A:332:LEU:HB2	1.80	0.63
1:A:15:VAL:HG13	1:A:27:VAL:HB	1.80	0.63
1:A:409:GLN:HA	1:A:409:GLN:NE2	2.14	0.63
1:A:22:LEU:HD12	1:A:26:ARG:CZ	2.29	0.63
2:D:97:VAL:HG23	2:D:103:ASN:HD22	1.62	0.63
1:A:81:ARG:HH22	1:A:102:GLY:H	1.47	0.62
2:D:16:LEU:CD2	2:D:31:LEU:HD21	2.29	0.62
2:D:48:ASN:OD1	2:D:190:GLN:HB2	1.98	0.62
1:A:370:ASN:C	1:A:370:ASN:HD22	2.07	0.62
1:A:133:ASN:HD21	1:A:392:THR:N	1.97	0.62
2:D:174:ARG:NH1	2:D:188:VAL:HG23	2.15	0.61
1:A:436:ASP:CG	1:A:437:ASN:H	2.09	0.61
1:A:522:VAL:HG11	1:A:526:GLN:NE2	2.14	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:345:GLU:O	2:D:352:VAL:HG12	2.00	0.61
1:A:182:ARG:HG2	1:A:182:ARG:HH11	1.64	0.61
1:A:195:GLU:HG2	1:A:197:TYR:CE2	2.34	0.61
1:A:343:VAL:HG13	1:A:343:VAL:O	2.01	0.61
2:D:58:LYS:HE3	2:D:58:LYS:O	1.99	0.61
2:D:97:VAL:HA	2:D:103:ASN:HB2	1.82	0.61
1:A:270:ASN:ND2	1:A:271:ARG:N	2.47	0.61
1:A:392:THR:HG23	1:A:402:GLU:HG2	1.82	0.61
1:A:557:MET:HE2	1:A:559:SER:HB2	1.83	0.61
1:A:537:LYS:C	1:A:537:LYS:HD3	2.26	0.61
1:A:77:VAL:HG21	1:A:121:TYR:CE1	2.37	0.60
1:A:131:VAL:HB	1:A:138:VAL:HG12	1.82	0.60
2:D:90:CYS:HB2	2:D:107:ILE:HD11	1.83	0.60
2:D:97:VAL:O	2:D:98:ASP:HB2	2.00	0.60
1:A:30:ARG:HB2	1:A:52:LEU:HD21	1.82	0.60
1:A:232:ASN:O	1:A:233:PRO:C	2.43	0.60
2:D:17:PHE:CD1	2:D:23:TRP:HA	2.37	0.60
2:D:31:LEU:N	2:D:31:LEU:HD22	2.17	0.60
2:D:143:ARG:HH21	2:D:143:ARG:HG3	1.67	0.60
2:D:174:ARG:HH11	2:D:188:VAL:HG23	1.67	0.59
1:A:133:ASN:HD22	1:A:390:VAL:HG12	1.67	0.59
1:A:243:THR:HG21	5:A:780:HOH:O	2.01	0.59
1:A:148:GLN:O	1:A:152:LEU:HD13	2.02	0.59
2:D:12:LEU:HD11	2:D:14:VAL:HG23	1.84	0.59
2:D:364:LEU:HD13	2:D:365:GLU:N	2.17	0.59
1:A:392:THR:OG1	1:A:402:GLU:HG2	2.02	0.59
1:A:320:THR:HB	1:A:324:ALA:HB2	1.83	0.59
2:D:73:SER:HB2	2:D:85:THR:HG23	1.84	0.59
2:D:40:MSE:HE3	2:D:86:ARG:HH21	1.68	0.58
2:D:112:ILE:O	2:D:116:GLU:HG3	2.03	0.58
2:D:173:VAL:HG13	2:D:177:ALA:HB2	1.84	0.58
1:A:50:GLU:CD	1:A:50:GLU:H	2.11	0.58
2:D:97:VAL:HA	2:D:103:ASN:CB	2.33	0.58
1:A:48:PRO:CG	1:A:51:LYS:HD3	2.33	0.58
2:D:41:ARG:HG2	2:D:343:ILE:HG12	1.86	0.58
2:D:360:ASN:H	2:D:360:ASN:ND2	2.02	0.58
1:A:328:THR:O	1:A:331:ILE:HG23	2.04	0.58
2:D:235:VAL:HG12	2:D:236:TRP:N	2.18	0.58
1:A:470:ARG:HG3	1:A:470:ARG:NH1	2.18	0.58
2:D:232:TYR:HD1	2:D:233:ASP:OD1	1.87	0.58
1:A:176:MET:HE3	1:A:180:LEU:CD2	2.33	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:41:ARG:HG2	2:D:343:ILE:HG23	1.85	0.58
2:D:80:ILE:H	2:D:80:ILE:CD1	2.17	0.58
2:D:160:ILE:HD12	2:D:160:ILE:O	2.04	0.58
2:D:20:TYR:O	2:D:23:TRP:HB3	2.04	0.57
1:A:200:ILE:HD13	1:A:236:ILE:HD12	1.86	0.57
1:A:25:GLY:HA3	1:A:73:VAL:HG12	1.87	0.57
1:A:132:LEU:O	1:A:390:VAL:HG11	2.05	0.57
1:A:425:SER:HB2	1:A:426:PRO:HD2	1.85	0.57
1:A:123:ARG:HD2	1:A:124:ARG:HH12	1.68	0.57
1:A:273:ARG:HG3	1:A:334:GLY:HA2	1.87	0.57
2:D:283:TYR:O	2:D:286:MSE:HB2	2.05	0.57
2:D:117:ASN:O	2:D:119:LYS:HG3	2.04	0.57
1:A:280:MET:HE1	1:A:295:MET:HE1	1.87	0.57
2:D:43:SER:HB3	2:D:47:ARG:HB3	1.87	0.57
2:D:91:LYS:HA	2:D:115:ILE:HD13	1.86	0.57
1:A:370:ASN:HD21	1:A:372:SER:HB3	1.69	0.56
2:D:241:ASN:C	2:D:241:ASN:HD22	2.13	0.56
2:D:318:LEU:CD2	2:D:355:LEU:HD11	2.35	0.56
1:A:170:ASP:HA	1:A:270:ASN:HD21	1.69	0.56
2:D:112:ILE:HB	2:D:140:TYR:OH	2.05	0.56
2:D:185:PHE:CD2	2:D:297:ARG:HA	2.40	0.56
2:D:208:PRO:HB2	2:D:325:ILE:HB	1.87	0.56
2:D:342:VAL:HG23	2:D:354:HIS:O	2.06	0.56
1:A:220:VAL:O	1:A:224:GLN:HG3	2.04	0.56
1:A:82:VAL:CG2	1:A:96:VAL:HG13	2.35	0.56
1:A:307:ALA:CB	1:A:331:ILE:HD11	2.33	0.56
2:D:318:LEU:HD23	2:D:355:LEU:HD21	1.87	0.56
2:D:286:MSE:HA	2:D:286:MSE:CE	2.31	0.56
1:A:157:ASN:HA	1:A:375:LEU:HG	1.88	0.56
1:A:60:PRO:C	1:A:62:THR:H	2.13	0.56
1:A:315:LEU:O	1:A:319:GLN:HG2	2.06	0.56
2:D:211:MSE:HE1	2:D:312:THR:HB	1.87	0.56
2:D:366:THR:O	2:D:369:PRO:HD3	2.06	0.56
1:A:236:ILE:HG23	1:A:240:GLU:HB2	1.87	0.55
1:A:227:ARG:NH2	1:A:229:ILE:HD13	2.22	0.55
2:D:57:SER:HB3	2:D:102:ASP:OD2	2.06	0.55
1:A:424:VAL:HA	1:A:429:ARG:O	2.07	0.55
2:D:133:LYS:CG	2:D:150:ILE:HG21	2.36	0.55
1:A:179:MET:HE3	1:A:275:MET:HE2	1.88	0.55
2:D:17:PHE:CE1	2:D:28:PHE:HB3	2.41	0.55
2:D:237:LEU:HD21	2:D:309:LYS:HG2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:429:ARG:HG2	1:A:429:ARG:HH11	1.71	0.55
1:A:58:ILE:HG12	1:A:97:LEU:CD2	2.36	0.55
2:D:30:GLU:C	2:D:31:LEU:HD22	2.32	0.55
1:A:66:MET:HE3	1:A:66:MET:HA	1.89	0.55
1:A:206:ILE:HG12	1:A:221:LEU:HD23	1.89	0.55
1:A:225:VAL:HG21	1:A:240:GLU:OE1	2.06	0.55
2:D:254:TYR:HB2	2:D:276:VAL:HG12	1.89	0.55
1:A:213:ASP:O	1:A:217:ILE:HG13	2.08	0.54
1:A:270:ASN:HD22	1:A:271:ARG:H	1.52	0.54
1:A:221:LEU:HD12	1:A:229:ILE:CG1	2.32	0.54
1:A:392:THR:CG2	1:A:402:GLU:HG2	2.38	0.54
1:A:416:HIS:CD2	1:A:417:PRO:HD2	2.42	0.54
2:D:189:TRP:CD1	2:D:320:PRO:HG2	2.43	0.54
2:D:279:ARG:CZ	2:D:286:MSE:HG2	2.38	0.54
1:A:370:ASN:ND2	1:A:372:SER:H	2.05	0.53
2:D:97:VAL:HG23	2:D:103:ASN:ND2	2.23	0.53
2:D:184:LYS:HE3	2:D:367:ILE:HD12	1.89	0.53
1:A:34:LEU:HD12	1:A:34:LEU:H	1.73	0.53
1:A:370:ASN:C	1:A:370:ASN:ND2	2.66	0.53
1:A:58:ILE:HG12	1:A:97:LEU:HD23	1.90	0.53
1:A:486:GLU:HG2	5:A:801:HOH:O	2.09	0.53
1:A:38:GLN:CD	1:A:38:GLN:H	2.16	0.53
2:D:237:LEU:H	2:D:237:LEU:CD2	2.21	0.53
2:D:245:ARG:HD3	2:D:305:ASP:OD1	2.08	0.53
2:D:11:ASN:HD22	2:D:75:ALA:HB3	1.74	0.53
2:D:43:SER:CB	2:D:47:ARG:HB3	2.38	0.53
2:D:157:VAL:HG13	2:D:172:TYR:HE2	1.73	0.53
1:A:265:VAL:O	1:A:269:VAL:HG22	2.09	0.53
1:A:256:ILE:HG23	1:A:257:LYS:N	2.22	0.53
1:A:373:ARG:HA	1:A:373:ARG:HE	1.74	0.53
2:D:96:SER:HG	2:D:104:ILE:HG23	1.74	0.53
2:D:178:LEU:HD23	2:D:293:GLU:HB3	1.90	0.53
2:D:217:SER:HB2	5:D:463:HOH:O	2.09	0.53
2:D:16:LEU:HA	2:D:69:ILE:O	2.09	0.52
2:D:61:HIS:CD2	2:D:95:VAL:HG11	2.44	0.52
2:D:250:ASN:ND2	2:D:272:GLU:HB2	2.23	0.52
1:A:81:ARG:HH22	1:A:102:GLY:N	2.05	0.52
2:D:367:ILE:HG13	5:D:431:HOH:O	2.09	0.52
1:A:22:LEU:HB3	1:A:24:LEU:HD22	1.91	0.52
1:A:263:GLY:N	1:A:264:PRO:HD2	2.25	0.52
1:A:59:GLN:HB3	1:A:60:PRO:HD2	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:MET:HG3	1:A:180:LEU:CD2	2.39	0.52
2:D:87:ILE:HD12	2:D:87:ILE:N	2.21	0.52
1:A:82:VAL:HG21	1:A:96:VAL:CG1	2.40	0.52
1:A:60:PRO:C	1:A:62:THR:N	2.67	0.52
1:A:254:ARG:HH11	1:A:254:ARG:HG2	1.75	0.52
2:D:330:LYS:HE3	2:D:330:LYS:CA	2.34	0.52
2:D:346:LEU:N	2:D:346:LEU:HD23	2.24	0.52
1:A:127:ASN:HD22	1:A:127:ASN:C	2.18	0.52
1:A:292:PHE:HB3	1:A:295:MET:HB3	1.91	0.52
2:D:374:ASN:C	2:D:376:PHE:H	2.17	0.52
1:A:31:VAL:CB	1:A:34:LEU:HD11	2.36	0.52
1:A:284:MET:HE1	1:A:292:PHE:CE1	2.42	0.51
1:A:402:GLU:OE1	1:A:411:ARG:NH1	2.43	0.51
1:A:557:MET:HE3	1:A:558:SER:H	1.75	0.51
2:D:12:LEU:HD11	2:D:14:VAL:CG2	2.40	0.51
2:D:38:LEU:HD13	2:D:52:LEU:HG	1.93	0.51
1:A:123:ARG:HD2	1:A:124:ARG:NH1	2.25	0.51
1:A:321:LYS:O	1:A:322:GLY:C	2.54	0.51
2:D:157:VAL:CG1	2:D:172:TYR:HE2	2.24	0.51
1:A:176:MET:O	1:A:180:LEU:CD2	2.57	0.51
2:D:53:GLN:HB3	2:D:106:ALA:HB2	1.93	0.51
1:A:85:HIS:CD2	1:A:97:LEU:HD11	2.46	0.51
1:A:199:THR:HG22	1:A:200:ILE:N	2.26	0.51
2:D:346:LEU:HB3	2:D:351:SER:CB	2.41	0.50
1:A:77:VAL:CG2	1:A:121:TYR:CE1	2.94	0.50
1:A:238:MET:O	1:A:242:THR:HG22	2.11	0.50
1:A:378:GLU:OE1	1:A:433:LYS:HG3	2.11	0.50
1:A:395:THR:OG1	1:A:399:HIS:HB2	2.11	0.50
2:D:136:LEU:HD13	2:D:150:ILE:HD11	1.94	0.50
1:A:522:VAL:HG11	1:A:526:GLN:HE21	1.75	0.50
2:D:154:ASN:CG	2:D:180:VAL:HG21	2.37	0.50
2:D:211:MSE:HB2	2:D:326:PHE:CE1	2.47	0.50
2:D:312:THR:CG2	2:D:353:THR:HB	2.42	0.50
1:A:332:LEU:O	1:A:332:LEU:HD22	2.11	0.50
1:A:557:MET:HE3	1:A:558:SER:N	2.27	0.50
2:D:80:ILE:HD13	2:D:80:ILE:N	2.26	0.50
2:D:174:ARG:HG3	2:D:188:VAL:HG22	1.94	0.50
2:D:213:VAL:HG22	2:D:213:VAL:O	2.12	0.50
2:D:268:THR:HG23	2:D:299:GLN:NE2	2.27	0.50
1:A:200:ILE:HG12	1:A:201:GLY:N	2.23	0.49
1:A:154:THR:H	1:A:182:ARG:NH2	2.09	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:133:LYS:HE3	2:D:152:ALA:HB3	1.94	0.49
1:A:102:GLY:O	1:A:128:ASP:HB2	2.12	0.49
1:A:60:PRO:HG2	1:A:62:THR:HG22	1.94	0.49
1:A:423:GLU:OE1	1:A:429:ARG:NH2	2.46	0.49
2:D:345:GLU:HB3	2:D:352:VAL:CG1	2.43	0.49
1:A:207:MET:SD	1:A:211:VAL:HG21	2.53	0.49
2:D:11:ASN:H	2:D:74:VAL:CG2	2.19	0.49
1:A:61:ILE:C	1:A:61:ILE:HD12	2.37	0.49
1:A:553:LYS:HB2	1:A:553:LYS:NZ	2.27	0.49
1:A:225:VAL:HG11	1:A:229:ILE:HD11	1.95	0.49
1:A:270:ASN:O	1:A:274:GLN:HB2	2.12	0.49
1:A:315:LEU:O	1:A:315:LEU:HD13	2.13	0.49
2:D:167:GLU:CD	2:D:167:GLU:H	2.21	0.49
1:A:25:GLY:CA	1:A:73:VAL:HG12	2.43	0.48
1:A:202:ILE:O	1:A:202:ILE:HG22	2.13	0.48
1:A:256:ILE:HA	1:A:286:VAL:CG2	2.36	0.48
1:A:499:LYS:NZ	5:A:599:HOH:O	2.45	0.48
2:D:11:ASN:ND2	2:D:80:ILE:HG22	2.26	0.48
1:A:18:ARG:NH1	1:A:77:VAL:HG23	2.28	0.48
1:A:170:ASP:C	1:A:172:PRO:HD3	2.38	0.48
2:D:185:PHE:CE2	2:D:297:ARG:HA	2.48	0.48
1:A:58:ILE:HG13	1:A:97:LEU:HA	1.95	0.48
1:A:172:PRO:HB3	1:A:336:LEU:HD22	1.94	0.48
1:A:280:MET:CE	1:A:295:MET:HE1	2.43	0.48
2:D:215:GLU:H	2:D:329:SER:HB2	1.77	0.48
1:A:182:ARG:HG2	1:A:182:ARG:NH1	2.29	0.48
2:D:173:VAL:HG13	2:D:177:ALA:CB	2.43	0.48
2:D:179:ALA:HB1	2:D:182:SER:HB3	1.95	0.48
2:D:313:ILE:HD12	2:D:313:ILE:N	2.27	0.48
1:A:133:ASN:ND2	1:A:390:VAL:HG12	2.28	0.48
1:A:295:MET:HE2	1:A:331:ILE:HD12	1.95	0.48
2:D:17:PHE:CD1	2:D:28:PHE:HB3	2.48	0.48
1:A:127:ASN:ND2	1:A:129:THR:H	2.11	0.48
1:A:162:PRO:HG2	1:A:165:GLU:HG3	1.96	0.48
2:D:158:PRO:HD3	2:D:286:MSE:HE1	1.95	0.48
1:A:284:MET:HE2	1:A:288:GLY:C	2.39	0.48
2:D:12:LEU:HA	2:D:74:VAL:HB	1.96	0.48
2:D:58:LYS:HB3	2:D:60:ILE:HG13	1.96	0.48
1:A:337:GLU:OE1	1:A:343:VAL:HG12	2.14	0.48
2:D:248:MSE:HG2	2:D:298:LEU:O	2.13	0.48
1:A:18:ARG:NH1	1:A:78:GLU:OE2	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:SER:OG	1:A:178:GLU:HG3	2.14	0.47
2:D:40:MSE:HG2	2:D:88:TYR:OH	2.14	0.47
1:A:236:ILE:HG23	1:A:240:GLU:CB	2.45	0.47
1:A:22:LEU:HD12	1:A:26:ARG:NH2	2.30	0.47
2:D:75:ALA:HB2	2:D:83:VAL:HA	1.95	0.47
1:A:209:GLN:HB3	1:A:211:VAL:HG23	1.97	0.47
2:D:346:LEU:HB2	2:D:350:ASN:O	2.15	0.47
2:D:144:THR:O	2:D:147:THR:HG22	2.15	0.47
2:D:369:PRO:O	2:D:370:VAL:C	2.57	0.47
1:A:375:LEU:HD22	1:A:455:LYS:HG2	1.97	0.47
2:D:119:LYS:NZ	5:D:392:HOH:O	2.47	0.47
2:D:204:ILE:C	2:D:206:GLN:H	2.23	0.47
1:A:11:PHE:CD1	1:A:11:PHE:C	2.93	0.47
1:A:28:ARG:HD3	1:A:49:THR:O	2.15	0.47
1:A:239:GLU:O	1:A:242:THR:HG23	2.15	0.47
1:A:362:ASP:N	1:A:363:PRO:CD	2.78	0.47
2:D:5:GLN:HG3	2:D:9:TYR:CD1	2.50	0.47
1:A:498:VAL:HG11	1:A:502:VAL:HG23	1.97	0.47
1:A:538:VAL:HG11	1:A:542:VAL:HG12	1.96	0.47
1:A:7:ASN:HD22	1:A:7:ASN:HA	1.48	0.47
1:A:58:ILE:C	1:A:58:ILE:HD12	2.40	0.47
1:A:214:MET:HG3	1:A:218:ASN:ND2	2.30	0.47
2:D:16:LEU:H	2:D:16:LEU:HD23	1.79	0.47
2:D:237:LEU:HD22	2:D:237:LEU:N	2.27	0.47
1:A:81:ARG:HB3	1:A:100:TYR:CZ	2.50	0.46
1:A:564:GLN:HG3	1:A:566:THR:HG22	1.96	0.46
1:A:252:MET:CE	1:A:253:GLN:HG3	2.45	0.46
1:A:323:ARG:HA	1:A:326:ARG:HH11	1.80	0.46
1:A:92:THR:HB	2:D:282:ALA:CB	2.45	0.46
2:D:260:ASP:OD1	2:D:289:ARG:NH2	2.48	0.46
1:A:207:MET:O	1:A:207:MET:HG3	2.14	0.46
2:D:318:LEU:HB3	2:D:355:LEU:HD21	1.98	0.46
1:A:92:THR:HB	2:D:282:ALA:HB3	1.96	0.46
2:D:114:SER:HA	2:D:167:GLU:OE2	2.16	0.46
2:D:254:TYR:HE2	2:D:274:SER:HB2	1.81	0.46
2:D:125:PHE:CE2	2:D:134:GLU:HG3	2.50	0.46
1:A:383:TYR:O	1:A:384:LYS:HG3	2.15	0.46
2:D:41:ARG:CG	2:D:343:ILE:HG23	2.46	0.46
2:D:91:LYS:O	2:D:91:LYS:HD3	2.16	0.46
2:D:153:ILE:HG12	2:D:197:MSE:HB3	1.98	0.46
1:A:416:HIS:HD2	1:A:418:THR:N	2.11	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:57:SER:C	2:D:59:ASN:N	2.74	0.46
1:A:214:MET:HG3	1:A:218:ASN:HD21	1.82	0.45
1:A:300:LEU:C	1:A:302:GLY:H	2.24	0.45
2:D:54:PHE:CE1	2:D:105:ILE:HD11	2.50	0.45
2:D:122:ARG:HG3	2:D:123:PRO:O	2.17	0.45
1:A:9:ASN:HD22	1:A:9:ASN:HA	1.53	0.45
1:A:85:HIS:NE2	1:A:97:LEU:HD11	2.32	0.45
2:D:56:ASP:HA	5:D:413:HOH:O	2.16	0.45
1:A:78:GLU:H	1:A:78:GLU:CD	2.24	0.45
1:A:456:THR:HG22	1:A:457:ASN:N	2.31	0.45
1:A:214:MET:CG	1:A:232:ASN:O	2.61	0.45
2:D:32:ALA:O	2:D:35:ILE:HG22	2.17	0.45
2:D:236:TRP:HH2	2:D:332:GLN:O	1.99	0.45
2:D:90:CYS:CB	2:D:107:ILE:HD11	2.47	0.45
2:D:143:ARG:HG3	2:D:143:ARG:NH2	2.30	0.45
2:D:345:GLU:HB3	2:D:352:VAL:HG13	1.99	0.45
1:A:121:TYR:HA	1:A:122:PRO:C	2.41	0.45
1:A:407:PRO:C	1:A:409:GLN:H	2.25	0.45
1:A:303:ASP:OD1	1:A:306:LYS:HB2	2.17	0.45
2:D:131:SER:O	2:D:135:MSE:HG3	2.17	0.45
2:D:363:LYS:HG2	5:D:435:HOH:O	2.16	0.45
1:A:337:GLU:HG2	1:A:342:GLU:HA	1.99	0.45
2:D:10:PRO:HG3	2:D:346:LEU:HD12	1.99	0.45
2:D:156:TYR:O	2:D:177:ALA:HA	2.17	0.45
2:D:19:SER:H	2:D:22:ALA:HB3	1.80	0.45
2:D:92:HIS:ND1	2:D:92:HIS:C	2.75	0.45
2:D:174:ARG:NH1	2:D:188:VAL:O	2.50	0.45
2:D:251:ALA:CB	2:D:273:ASN:HB2	2.46	0.45
1:A:41:GLN:CD	2:D:123:PRO:HD2	2.42	0.44
1:A:73:VAL:HG12	1:A:73:VAL:O	2.17	0.44
2:D:91:LYS:CA	2:D:115:ILE:HD13	2.47	0.44
2:D:250:ASN:O	2:D:273:ASN:HB2	2.17	0.44
2:D:319:THR:O	2:D:322:VAL:HG12	2.16	0.44
1:A:554:MET:HE3	1:A:554:MET:HB3	1.89	0.44
2:D:198:MSE:HE1	2:D:206:GLN:HE22	1.82	0.44
2:D:199:ASP:OD2	2:D:199:ASP:C	2.60	0.44
1:A:17:ASP:HB3	1:A:28:ARG:HB2	1.98	0.44
2:D:244:LYS:O	2:D:247:PRO:HD2	2.17	0.44
1:A:285:GLY:C	1:A:287:GLY:N	2.75	0.44
2:D:279:ARG:HG3	2:D:289:ARG:O	2.17	0.44
2:D:66:GLY:O	2:D:90:CYS:HB3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:295:ALA:O	2:D:299:GLN:HG3	2.18	0.44
2:D:296:ILE:HD11	2:D:369:PRO:HD2	1.98	0.44
2:D:328:ASP:HB2	2:D:333:PHE:CE2	2.53	0.44
1:A:368:ILE:O	1:A:368:ILE:HG13	2.16	0.44
1:A:416:HIS:HD2	1:A:418:THR:CB	2.13	0.44
1:A:31:VAL:O	1:A:35:HIS:HB2	2.18	0.44
2:D:43:SER:HA	2:D:341:GLU:HB3	2.00	0.44
2:D:189:TRP:NE1	2:D:320:PRO:HG2	2.33	0.44
1:A:99:THR:HG22	1:A:100:TYR:N	2.33	0.44
1:A:200:ILE:HG21	1:A:236:ILE:HD12	1.99	0.44
1:A:370:ASN:HD21	1:A:372:SER:HB2	1.79	0.44
1:A:418:THR:HG22	1:A:419:GLY:N	2.33	0.44
1:A:436:ASP:CG	1:A:437:ASN:N	2.74	0.44
2:D:213:VAL:O	2:D:213:VAL:HG13	2.16	0.44
2:D:21:ASP:C	2:D:23:TRP:N	2.73	0.44
1:A:253:GLN:HE21	1:A:253:GLN:HB2	1.61	0.43
1:A:409:GLN:NE2	1:A:409:GLN:CA	2.77	0.43
2:D:20:TYR:CZ	2:D:143:ARG:HA	2.54	0.43
2:D:35:ILE:O	2:D:348:ASN:ND2	2.51	0.43
2:D:262:SER:C	2:D:264:PRO:HD3	2.43	0.43
1:A:320:THR:HG22	1:A:323:ARG:HB3	1.99	0.43
1:A:541:THR:HG23	5:A:848:HOH:O	2.18	0.43
1:A:242:THR:HG23	1:A:243:THR:H	1.83	0.43
2:D:35:ILE:HD11	2:D:38:LEU:HB2	2.00	0.43
2:D:95:VAL:HG13	2:D:105:ILE:CG2	2.48	0.43
2:D:125:PHE:HE2	2:D:134:GLU:HG3	1.82	0.43
2:D:235:VAL:HG12	2:D:236:TRP:H	1.83	0.43
2:D:246:ASP:HB2	2:D:247:PRO:HD3	1.99	0.43
1:A:439:TYR:HB3	1:A:441:ILE:CD1	2.49	0.43
2:D:114:SER:HA	2:D:167:GLU:CD	2.44	0.43
2:D:60:ILE:HG22	2:D:64:MSE:HG3	2.00	0.43
2:D:174:ARG:CG	2:D:188:VAL:HG22	2.48	0.43
1:A:468:ASP:CG	1:A:469:ASN:H	2.27	0.43
2:D:244:LYS:C	2:D:247:PRO:HD2	2.43	0.43
1:A:82:VAL:HG22	1:A:83:TYR:N	2.34	0.43
1:A:48:PRO:HB2	1:A:50:GLU:OE2	2.19	0.43
1:A:299:MET:HG3	1:A:304:TRP:HZ3	1.84	0.43
2:D:47:ARG:CB	2:D:47:ARG:NH1	2.82	0.43
1:A:43:ASP:O	2:D:122:ARG:NE	2.52	0.42
1:A:270:ASN:O	1:A:274:GLN:N	2.52	0.42
2:D:28:PHE:HD1	2:D:29:VAL:N	2.17	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:236:TRP:CD1	2:D:308:ALA:HB2	2.54	0.42
1:A:332:LEU:HD13	1:A:333:THR:HG23	1.99	0.42
2:D:245:ARG:HD3	2:D:305:ASP:CG	2.44	0.42
1:A:239:GLU:CD	1:A:239:GLU:H	2.28	0.42
1:A:216:GLN:O	1:A:219:LYS:HB3	2.20	0.42
1:A:242:THR:O	1:A:245:PHE:N	2.51	0.42
2:D:17:PHE:CD2	2:D:17:PHE:N	2.88	0.42
2:D:143:ARG:HB3	2:D:146:LEU:HD12	2.01	0.42
2:D:184:LYS:HE3	2:D:367:ILE:CG2	2.49	0.42
2:D:212:ILE:HG13	2:D:230:LEU:HD13	2.01	0.42
2:D:306:GLY:O	2:D:358:PHE:HA	2.20	0.42
1:A:36:PRO:HB2	1:A:38:GLN:OE1	2.20	0.42
1:A:209:GLN:HE21	1:A:209:GLN:HB2	1.64	0.42
1:A:213:ASP:OD2	1:A:216:GLN:HG3	2.20	0.42
2:D:292:TYR:CZ	2:D:296:ILE:HG13	2.54	0.42
2:D:324:ILE:HB	2:D:357:MSE:CE	2.50	0.42
2:D:136:LEU:CD1	2:D:150:ILE:HD11	2.49	0.42
1:A:214:MET:HG2	1:A:232:ASN:C	2.43	0.42
1:A:278:GLU:O	1:A:281:ALA:HB3	2.19	0.42
1:A:296:LEU:HD12	1:A:296:LEU:HA	1.89	0.42
2:D:209:TYR:OH	2:D:229:PRO:HG2	2.19	0.42
2:D:318:LEU:HD23	2:D:355:LEU:HD11	2.02	0.42
1:A:255:ASP:HB3	1:A:286:VAL:HG11	2.02	0.42
1:A:383:TYR:CD1	1:A:383:TYR:C	2.98	0.42
2:D:56:ASP:OD2	2:D:60:ILE:HG13	2.20	0.42
2:D:97:VAL:HG22	2:D:98:ASP:N	2.35	0.42
2:D:241:ASN:C	2:D:241:ASN:ND2	2.74	0.41
1:A:286:VAL:HG12	5:A:692:HOH:O	2.21	0.41
1:A:326:ARG:O	1:A:330:ILE:HG12	2.21	0.41
1:A:402:GLU:CD	1:A:411:ARG:HH12	2.28	0.41
1:A:416:HIS:CG	1:A:417:PRO:HD2	2.54	0.41
2:D:349:ASN:OD1	2:D:350:ASN:HB2	2.21	0.41
1:A:184:GLU:HA	5:A:749:HOH:O	2.19	0.41
2:D:23:TRP:HZ3	2:D:145:LEU:CD1	2.33	0.41
2:D:47:ARG:HB3	2:D:47:ARG:HH11	1.85	0.41
2:D:57:SER:CB	2:D:102:ASP:OD2	2.69	0.41
1:A:21:PRO:HD3	1:A:54:TRP:CZ2	2.55	0.41
2:D:11:ASN:C	2:D:74:VAL:HG23	2.45	0.41
2:D:128:ALA:O	2:D:132:ILE:HG13	2.20	0.41
2:D:160:ILE:HD12	2:D:160:ILE:C	2.46	0.41
1:A:242:THR:O	1:A:245:PHE:HB3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:265:VAL:HG22	1:A:300:LEU:HD23	2.03	0.41
1:A:323:ARG:HA	1:A:326:ARG:NH1	2.36	0.41
1:A:456:THR:CG2	1:A:457:ASN:N	2.84	0.41
2:D:154:ASN:O	2:D:180:VAL:HG23	2.20	0.41
1:A:232:ASN:HB3	1:A:233:PRO:HD3	2.03	0.41
1:A:385:GLY:HA2	1:A:406:THR:OG1	2.21	0.41
2:D:34:THR:HB	2:D:55:TYR:O	2.21	0.41
2:D:133:LYS:HE3	2:D:152:ALA:CB	2.51	0.41
2:D:208:PRO:O	2:D:210:PRO:CD	2.67	0.41
1:A:26:ARG:HD3	2:D:259:LEU:HD21	2.02	0.40
1:A:176:MET:HG3	1:A:180:LEU:HD21	2.02	0.40
1:A:323:ARG:CA	1:A:326:ARG:HH11	2.34	0.40
1:A:421:TYR:CE1	1:A:433:LYS:HB2	2.56	0.40
2:D:361:ALA:C	2:D:362:THR:HG22	2.46	0.40
2:D:113:HIS:HB3	2:D:139:ILE:HG21	2.03	0.40
2:D:179:ALA:CB	2:D:182:SER:HB3	2.51	0.40
2:D:235:VAL:O	2:D:308:ALA:HB1	2.21	0.40
1:A:252:MET:CE	1:A:278:GLU:HB3	2.51	0.40
2:D:12:LEU:CD1	2:D:14:VAL:HG23	2.51	0.40
2:D:235:VAL:CG1	2:D:236:TRP:N	2.84	0.40
2:D:254:TYR:CE2	2:D:274:SER:HB2	2.56	0.40
1:A:137:GLU:O	1:A:141:ASP:N	2.48	0.40
2:D:157:VAL:HG13	2:D:172:TYR:CE2	2.53	0.40
2:D:368:ASP:N	2:D:369:PRO:CD	2.84	0.40
2:D:28:PHE:CD1	2:D:28:PHE:C	3.00	0.40
2:D:41:ARG:NH2	2:D:341:GLU:OE2	2.54	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:586:PO4:P	4:A:586:PO4:O2[2_555]	1.48	0.72
1:A:584:HIS:OXT	1:A:584:HIS:OXT[4_555]	1.87	0.33

5.3 Torsion angles

5.3.1 Protein backbone

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	551/584 (94%)	512 (93%)	31 (6%)	8 (2%)	8	28
2	D	360/391 (92%)	295 (82%)	52 (14%)	13 (4%)	2	11
All	All	911/975 (93%)	807 (89%)	83 (9%)	21 (2%)	5	19

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	575	GLY
2	D	77	ALA
2	D	213	VAL
1	A	232	ASN
2	D	98	ASP
2	D	99	SER
2	D	362	THR
2	D	367	ILE
2	D	370	VAL
1	A	301	ALA
2	D	243	HIS
1	A	110	ARG
1	A	384	LYS
2	D	100	LYS
2	D	124	PHE
2	D	92	HIS
1	A	322	GLY
2	D	313	ILE
1	A	200	ILE
1	A	202	ILE
2	D	74	VAL

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	472/494 (96%)	427 (90%)	45 (10%)	8	26
2	D	324/333 (97%)	305 (94%)	19 (6%)	18	48
All	All	796/827 (96%)	732 (92%)	64 (8%)	11	34

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	9	ASN
1	A	18	ARG
1	A	24	LEU
1	A	34	LEU
1	A	45	MET
1	A	50	GLU
1	A	52	LEU
1	A	59	GLN
1	A	62	THR
1	A	66	MET
1	A	78	GLU
1	A	87	LEU
1	A	89	LYS
1	A	129	THR
1	A	152	LEU
1	A	184	GLU
1	A	195	GLU
1	A	209	GLN
1	A	242	THR
1	A	296	LEU
1	A	300	LEU
1	A	327	VAL
1	A	331	ILE
1	A	332	LEU
1	A	335	ASN
1	A	375	LEU
1	A	388	PRO
1	A	393	MET
1	A	397	SER
1	A	411	ARG
1	A	418	THR
1	A	429	ARG
1	A	433	LYS
1	A	442	THR

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Mol	Chain	Res	Type
1	A	457	ASN
1	A	482	VAL
1	A	486	GLU
1	A	495	THR
1	A	519	THR
1	A	522	VAL
1	A	542	VAL
1	A	550	TRP
1	A	553	LYS
1	A	566	THR
2	D	31	LEU
2	D	44	LEU
2	D	58	LYS
2	D	68	GLU
2	D	80	ILE
2	D	87	ILE
2	D	91	LYS
2	D	183	ASP
2	D	236	TRP
2	D	237	LEU
2	D	267	THR
2	D	289	ARG
2	D	293	GLU
2	D	296	ILE
2	D	325	ILE
2	D	330	LYS
2	D	350	ASN
2	D	357	MSE
2	D	360	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (39) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	9	ASN
1	A	59	GLN
1	A	85	HIS
1	A	93	ASN
1	A	127	ASN
1	A	133	ASN
1	A	173	ASN
1	A	209	GLN

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Mol	Chain	Res	Type
1	A	218	ASN
1	A	253	GLN
1	A	270	ASN
1	A	318	GLN
1	A	335	ASN
1	A	370	ASN
1	A	409	GLN
1	A	416	HIS
1	A	443	ASN
1	A	457	ASN
1	A	526	GLN
1	A	533	ASN
2	D	11	ASN
2	D	26	ASN
2	D	53	GLN
2	D	61	HIS
2	D	67	ASN
2	D	103	ASN
2	D	154	ASN
2	D	206	GLN
2	D	241	ASN
2	D	273	ASN
2	D	303	GLN
2	D	315	ASN
2	D	331	ASN
2	D	332	GLN
2	D	350	ASN
2	D	354	HIS
2	D	360	ASN
2	D	374	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	$\# Z > 2$
4	PO4	A	586	-	0,2,4	-	-	0,1,6	-	-

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	586	PO4	0	1

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	557/584 (95%)	-0.29	2 (0%) 88 85	22, 45, 71, 103	0
2	D	349/391 (89%)	0.46	9 (2%) 57 48	49, 85, 120, 138	0
All	All	906/975 (92%)	-0.00	11 (1%) 76 69	22, 59, 111, 138	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	347	SER	2.6
2	D	90	CYS	2.5
2	D	97	VAL	2.4
2	D	376	PHE	2.4
2	D	77	ALA	2.4
2	D	45	TYR	2.3
1	A	269	VAL	2.1
2	D	99	SER	2.0
1	A	583	HIS	2.0
2	D	232	TYR	2.0
2	D	338	TYR	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
3	K	A	585	1/1	0.96	0.05	48,48,48,48	1
4	PO4	A	586	3/5	0.98	0.09	24,24,25,95	2

6.5 Other polymers [i](#)

There are no such residues in this entry.