



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 03:01 AM UTC

PDB ID : 1H4T / pdb_00001h4t
Title : Prolyl-tRNA synthetase from *Thermus thermophilus* complexed with L-proline
Authors : Yaremchuk, A.; Tukalo, M.; Cusack, S.
Deposited on : 2001-05-14
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

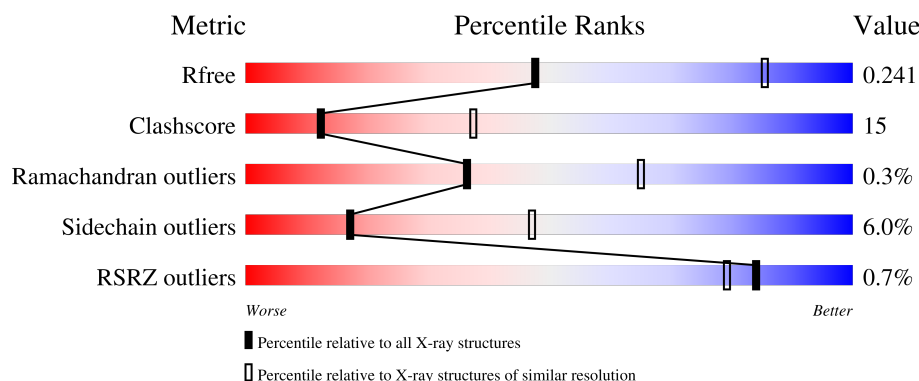
1 Overall quality at a glance

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	477	64%29% . .
1	B	477	% 61%31%5% .
1	C	477	69%25% . . .
1	D	477	% 68%25% . .

2 Entry composition [i](#)

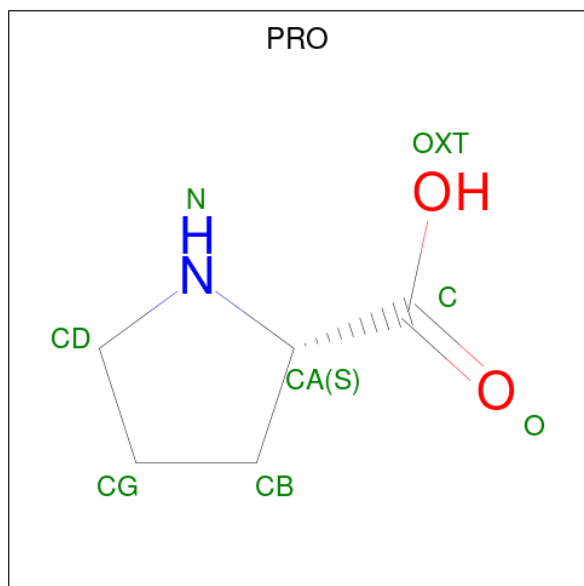
There are 4 unique types of molecules in this entry. The entry contains 15464 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PROLYL-TRNA SYNTHETASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	464	Total	C	N	O	S	42	0	0
			3747	2399	656	680	12			
1	B	464	Total	C	N	O	S	32	0	0
			3747	2399	656	680	12			
1	C	464	Total	C	N	O	S	30	0	0
			3747	2399	656	680	12			
1	D	464	Total	C	N	O	S	36	0	0
			3747	2399	656	680	12			

- Molecule 2 is PROLINE (CCD ID: PRO) (formula: C₅H₉NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			8	5	1	2		
2	B	1	Total	C	N	O	0	0
			8	5	1	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	N	O	0	0
			8	5	1	2		
2	D	1	Total	C	N	O	0	0
			8	5	1	2		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		
3	C	1	Total	Zn	0	0
			1	1		
3	D	1	Total	Zn	0	0
			1	1		

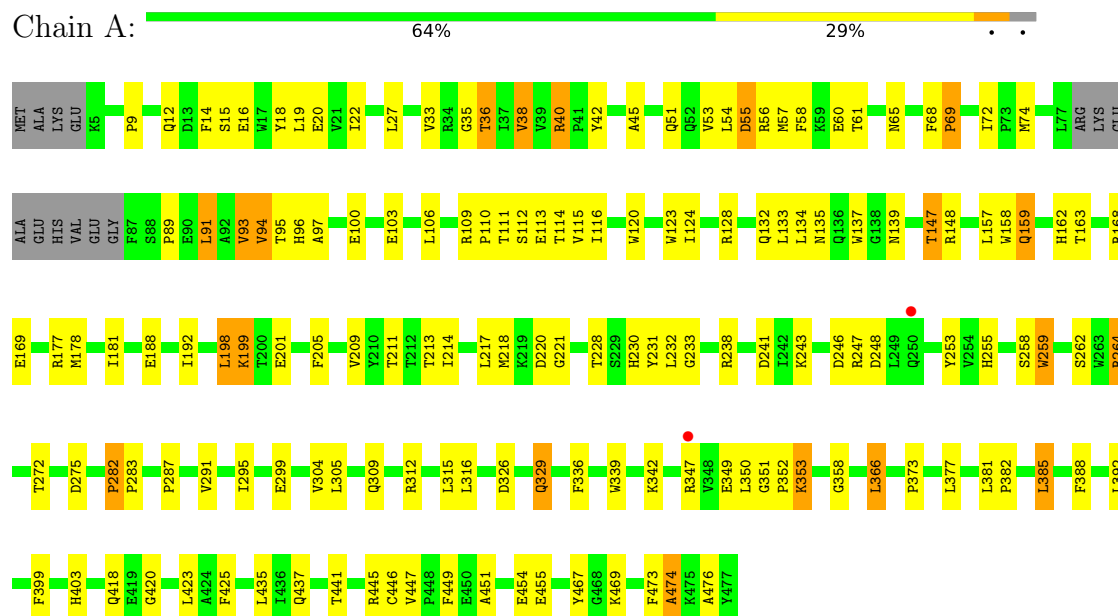
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	90	Total	O	0	0
			90	90		
4	B	113	Total	O	0	0
			113	113		
4	C	112	Total	O	0	0
			112	112		
4	D	125	Total	O	0	0
			125	125		

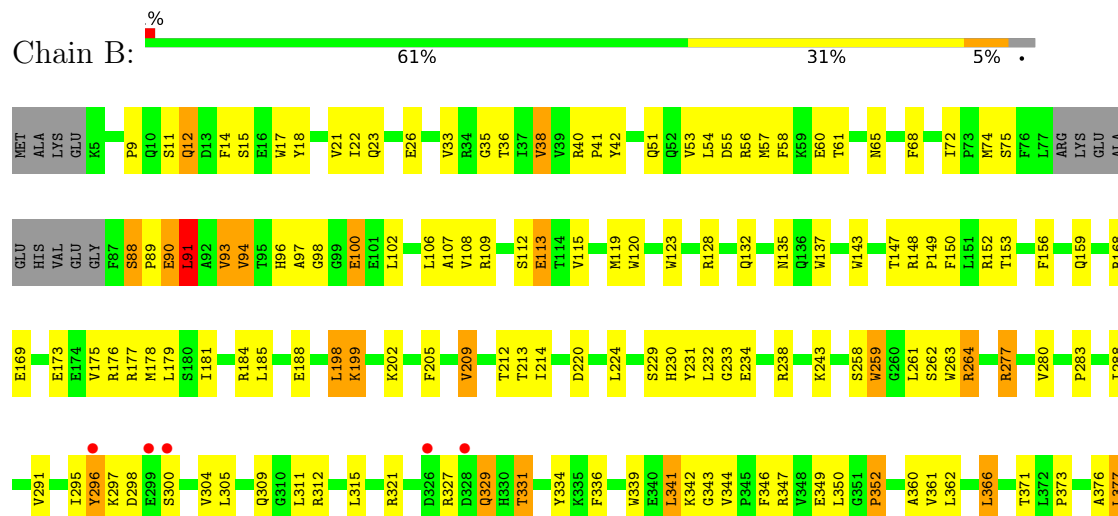
3 Residue-property plots [i](#)

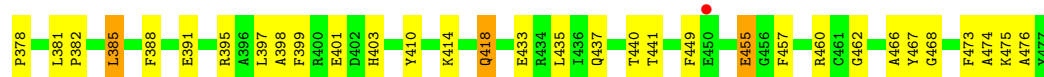
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

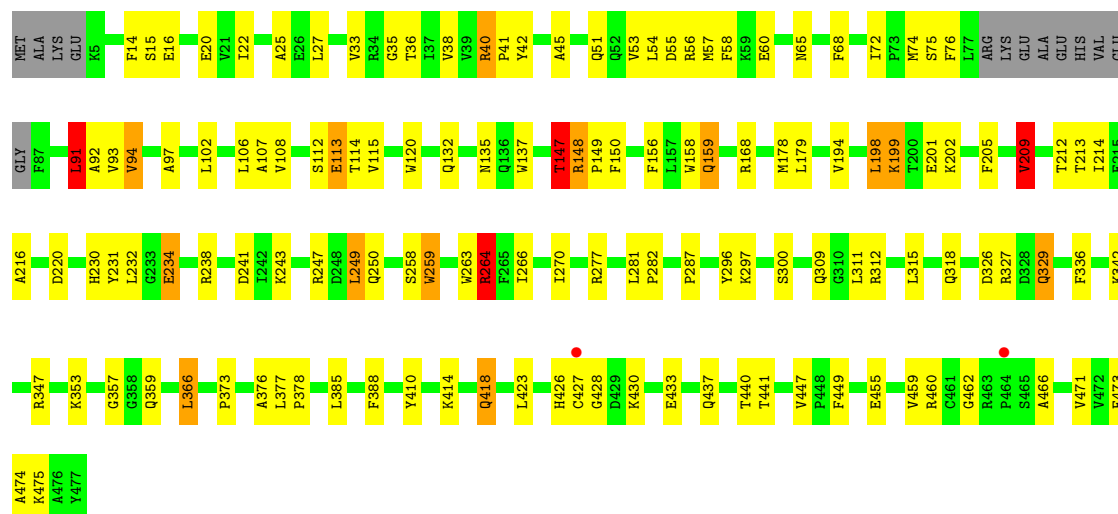


• Molecule 1: PROLYL-TRNA SYNTHETASE

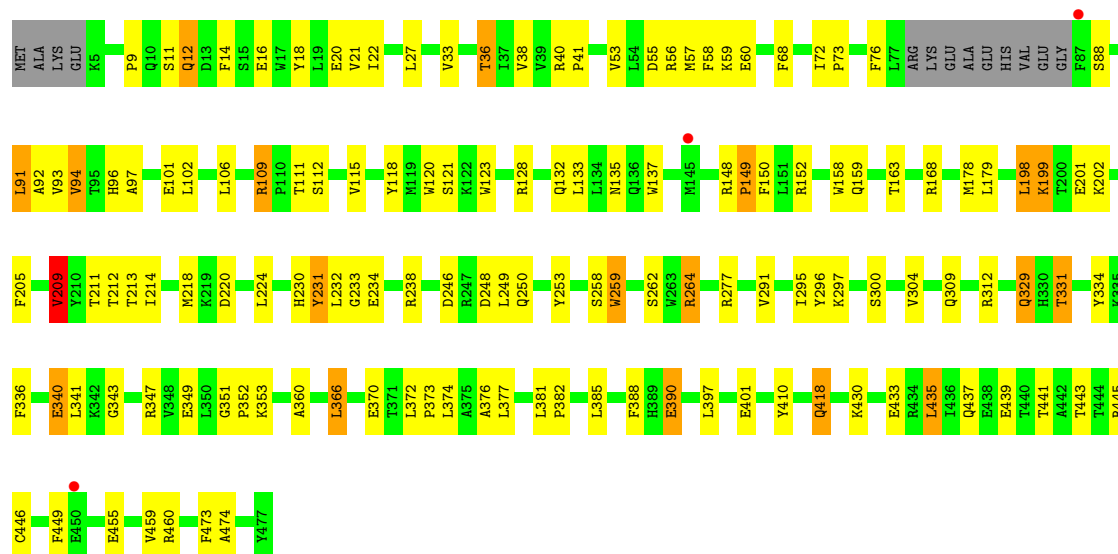




• Molecule 1: PROLYL-TRNA SYNTHETASE



• Molecule 1: PROLYL-TRNA SYNTHETASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	132.57Å 193.47Å 125.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.90 15.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.5 (15.00-2.90) 98.7 (15.00-2.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.98 (at 2.91Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.200 , 0.243 0.200 , 0.241	Depositor DCC
R_{free} test set	3538 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	39.6	Xtriage
Anisotropy	0.454	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 57.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	15464	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.88% 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:
ZN

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	0.39	0/3844	0.91	18/5207 (0.3%)
1	B	0.42	0/3844	0.95	17/5207 (0.3%)
1	C	0.43	0/3844	0.95	17/5207 (0.3%)
1	D	0.42	0/3844	0.95	16/5207 (0.3%)
All	All	0.42	0/15376	0.94	68/20828 (0.3%)

There are no bond length outliers.

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	209	VAL	N-CA-C	-11.15	100.00	110.82
1	A	209	VAL	N-CA-C	-8.38	103.12	110.74
1	D	112	SER	N-CA-C	7.97	121.80	112.72
1	C	91	LEU	N-CA-C	7.80	122.53	110.20
1	B	112	SER	N-CA-C	7.61	122.69	112.88
1	D	68	PHE	CA-C-N	7.34	127.38	119.90
1	D	68	PHE	C-N-CA	7.34	127.38	119.90
1	D	209	VAL	N-CA-C	-7.31	103.03	111.00
1	A	91	LEU	N-CA-C	6.93	120.54	110.28
1	B	264	ARG	N-CA-C	-6.92	104.10	112.54
1	C	264	ARG	N-CA-C	-6.86	103.83	111.71
1	C	148	ARG	CA-C-N	6.85	128.41	119.84
1	C	148	ARG	C-N-CA	6.85	128.41	119.84
1	B	209	VAL	N-CA-C	-6.75	103.64	111.00
1	D	91	LEU	N-CA-C	6.67	120.74	110.20
1	D	109	ARG	N-CA-C	6.57	119.04	109.71
1	C	68	PHE	CA-C-N	6.54	126.57	119.90
1	C	68	PHE	C-N-CA	6.54	126.57	119.90
1	A	112	SER	N-CA-C	6.49	119.85	112.97
1	C	475	LYS	N-CA-C	-6.43	98.79	109.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	264	ARG	N-CA-C	-6.28	104.88	112.54
1	D	93	VAL	N-CA-C	6.21	117.24	108.36
1	C	112	SER	N-CA-C	6.17	120.84	112.88
1	A	93	VAL	N-CA-C	6.06	117.30	108.58
1	D	340	GLU	N-CA-C	-5.99	104.67	111.14
1	B	109	ARG	N-CA-C	5.98	118.20	109.71
1	B	93	VAL	N-CA-C	5.96	116.88	108.36
1	C	418	GLN	N-CA-C	-5.95	105.85	113.23
1	B	91	LEU	N-CA-C	5.92	119.56	110.20
1	B	108	VAL	N-CA-C	-5.90	101.11	108.89
1	B	418	GLN	N-CA-C	-5.83	106.30	113.41
1	C	147	THR	N-CA-C	5.76	118.11	108.96
1	B	68	PHE	CA-C-N	5.72	125.73	119.89
1	B	68	PHE	C-N-CA	5.72	125.73	119.89
1	D	351	GLY	CA-C-N	5.69	126.96	119.84
1	D	351	GLY	C-N-CA	5.69	126.96	119.84
1	A	109	ARG	N-CA-C	5.68	116.56	109.57
1	D	264	ARG	N-CA-C	-5.67	105.46	112.38
1	D	92	ALA	N-CA-C	-5.58	98.12	107.99
1	D	199	LYS	N-CA-C	-5.56	102.29	110.52
1	B	199	LYS	N-CA-C	-5.56	102.25	110.48
1	C	199	LYS	N-CA-C	-5.55	102.31	110.52
1	A	351	GLY	CA-C-N	5.53	124.72	118.97
1	A	351	GLY	C-N-CA	5.53	124.72	118.97
1	C	281	LEU	N-CA-C	5.51	116.76	109.93
1	A	103	GLU	N-CA-C	-5.46	105.35	112.23
1	C	92	ALA	N-CA-C	-5.38	99.00	108.20
1	B	88	SER	CA-C-N	5.32	126.49	119.84
1	B	88	SER	C-N-CA	5.32	126.49	119.84
1	D	343	GLY	N-CA-C	5.28	122.08	114.10
1	B	475	LYS	N-CA-C	-5.28	101.36	109.76
1	B	476	ALA	N-CA-C	5.24	117.65	109.52
1	B	343	GLY	N-CA-C	5.20	121.95	114.10
1	C	282	PRO	CA-C-N	5.20	126.33	119.84
1	C	282	PRO	C-N-CA	5.20	126.33	119.84
1	A	114	THR	N-CA-C	5.19	116.62	110.97
1	A	188	GLU	N-CA-C	5.15	117.63	111.71
1	A	134	LEU	N-CA-C	5.12	117.91	109.72
1	A	476	ALA	N-CA-C	5.10	118.05	110.14
1	A	316	LEU	N-CA-C	-5.10	105.61	111.07
1	A	282	PRO	CA-C-N	5.09	126.21	119.84
1	A	282	PRO	C-N-CA	5.09	126.21	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	D	418	GLN	N-CA-C	-5.08	107.25	113.50
1	B	296	TYR	N-CA-C	5.08	115.85	108.14
1	C	108	VAL	N-CA-C	-5.08	102.14	108.84
1	D	231	TYR	N-CA-C	-5.08	100.46	108.73
1	A	69	PRO	N-CA-C	5.07	119.27	111.21
1	A	199	LYS	N-CA-C	-5.03	103.04	110.48

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	3747	0	3680	117	0
1	B	3747	0	3680	143	0
1	C	3747	0	3680	109	0
1	D	3747	0	3680	109	0
2	A	8	0	7	1	0
2	B	8	0	7	0	0
2	C	8	0	7	2	0
2	D	8	0	7	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	90	0	0	4	0
4	B	113	0	0	10	0
4	C	112	0	0	12	0
4	D	125	0	0	5	0
All	All	15464	0	14748	454	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 15.

All (454) 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:264:ARG:HG2	1:C:264:ARG:HH11	1.12	1.08
1:D:9:PRO:HG2	1:D:12:GLN:HB2	1.40	1.03
1:A:120:TRP:NE1	1:A:132:GLN:HG2	1.81	0.95
1:C:120:TRP:NE1	1:C:132:GLN:HG2	1.83	0.94
1:D:88:SER:HB3	1:D:91:LEU:HD23	1.50	0.91
1:D:120:TRP:NE1	1:D:132:GLN:HG2	1.85	0.91
1:B:120:TRP:NE1	1:B:132:GLN:HG2	1.87	0.90
1:C:33:VAL:O	1:C:36:THR:HG22	1.71	0.89
1:B:74:MET:HE2	1:B:107:ALA:HB2	1.57	0.87
1:C:347:ARG:HD3	4:C:2092:HOH:O	1.74	0.86
1:A:53:VAL:HG12	1:A:57:MET:HE2	1.58	0.85
1:B:33:VAL:O	1:B:36:THR:HG22	1.77	0.84
1:A:198:LEU:HD22	1:A:449:PHE:HE2	1.42	0.84
1:B:12:GLN:HG2	1:D:12:GLN:HG3	1.60	0.84
1:C:264:ARG:HH11	1:C:264:ARG:CG	1.91	0.83
1:D:59:LYS:HD2	4:D:2025:HOH:O	1.78	0.82
1:C:15:SER:HA	1:C:148:ARG:HH22	1.45	0.81
1:C:120:TRP:HE1	1:C:132:GLN:HG2	1.46	0.80
1:C:264:ARG:HG2	1:C:264:ARG:NH1	1.93	0.80
1:A:35:GLY:HA2	1:A:147:THR:HB	1.62	0.80
1:B:74:MET:HE1	1:B:91:LEU:HD23	1.62	0.79
1:A:199:LYS:HE2	1:A:213:THR:HG21	1.63	0.79
1:B:9:PRO:HG2	1:B:12:GLN:HB2	1.65	0.78
1:B:12:GLN:HG2	1:D:12:GLN:CG	2.15	0.77
1:A:33:VAL:O	1:A:36:THR:HB	1.85	0.77
1:C:199:LYS:HE2	1:C:213:THR:HG21	1.66	0.77
1:B:198:LEU:HD22	1:B:449:PHE:HE2	1.50	0.76
1:B:56:ARG:O	1:B:60:GLU:HG3	1.86	0.75
1:C:426:HIS:ND1	4:C:2105:HOH:O	2.19	0.74
1:C:168:ARG:HG3	1:C:231:TYR:OH	1.88	0.74
1:C:198:LEU:HD22	1:C:449:PHE:HE2	1.53	0.73
1:C:238:ARG:HD2	4:C:2070:HOH:O	1.88	0.73
1:C:309:GLN:HG2	1:C:312:ARG:HH22	1.53	0.73
1:C:56:ARG:O	1:C:60:GLU:HG3	1.89	0.72
1:B:53:VAL:HG12	1:B:57:MET:HE2	1.70	0.72
1:D:120:TRP:HE1	1:D:132:GLN:HG2	1.52	0.72
1:C:16:GLU:O	1:C:20:GLU:HG3	1.89	0.72
1:C:201:GLU:HB3	1:C:430:LYS:HE2	1.71	0.72
1:A:211:THR:HG23	1:A:230:HIS:ND1	2.04	0.72
1:C:15:SER:HA	1:C:148:ARG:NH2	2.05	0.71
1:A:198:LEU:HD22	1:A:449:PHE:CE2	2.26	0.70
1:A:15:SER:HA	1:A:148:ARG:HH22	1.56	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:THR:HG21	4:A:2005:HOH:O	1.91	0.70
1:B:96:HIS:HB3	4:B:2034:HOH:O	1.89	0.70
1:B:295:ILE:O	1:B:352:PRO:HD3	1.91	0.70
1:B:397:LEU:O	1:B:401:GLU:HG3	1.90	0.70
1:D:168:ARG:HG3	1:D:231:TYR:OH	1.92	0.70
1:D:460:ARG:HH11	1:D:460:ARG:HG3	1.57	0.69
1:D:36:THR:HG21	4:D:2010:HOH:O	1.91	0.69
1:D:94:VAL:HG13	1:D:106:LEU:HB2	1.73	0.69
1:A:349:GLU:HB3	4:A:2076:HOH:O	1.91	0.69
1:D:246:ASP:HB3	1:D:248:ASP:OD2	1.92	0.69
1:B:198:LEU:HD22	1:B:449:PHE:CE2	2.26	0.69
1:D:297:LYS:H	1:D:300:SER:HB2	1.55	0.69
1:A:19:LEU:HD21	1:A:148:ARG:NH1	2.08	0.69
1:B:414:LYS:O	1:B:418:GLN:HG2	1.93	0.68
1:A:120:TRP:HE1	1:A:132:GLN:HG2	1.54	0.68
1:C:309:GLN:HG2	1:C:312:ARG:NH2	2.08	0.68
1:D:198:LEU:HD22	1:D:449:PHE:HE2	1.58	0.68
1:C:94:VAL:HG13	1:C:106:LEU:HB2	1.75	0.67
1:A:241:ASP:OD2	1:A:243:LYS:HE3	1.95	0.67
1:B:72:ILE:HG12	1:B:115:VAL:HG21	1.75	0.67
1:D:459:VAL:HB	4:D:2115:HOH:O	1.95	0.67
1:B:168:ARG:HG3	1:B:231:TYR:OH	1.95	0.67
1:A:58:PHE:HE2	1:A:159:GLN:HG3	1.60	0.66
1:A:231:TYR:CZ	1:A:233:GLY:HA2	2.30	0.66
1:C:97:ALA:HB3	1:C:102:LEU:HD21	1.77	0.66
1:A:123:TRP:CZ2	1:B:40:ARG:HD3	2.30	0.66
1:D:16:GLU:O	1:D:20:GLU:HG3	1.94	0.66
1:D:94:VAL:CG1	1:D:106:LEU:HB2	2.25	0.66
1:A:283:PRO:HB3	1:A:385:LEU:HD12	1.76	0.66
1:A:16:GLU:O	1:A:20:GLU:HG3	1.94	0.65
1:C:35:GLY:HA2	1:C:147:THR:HB	1.78	0.65
1:A:38:VAL:HG11	1:B:119:MET:HE2	1.78	0.65
1:B:72:ILE:HG12	1:B:115:VAL:CG2	2.26	0.65
1:B:224:LEU:HD11	1:B:264:ARG:HE	1.62	0.65
1:C:336:PHE:CD1	1:C:347:ARG:HD2	2.33	0.64
1:B:291:VAL:HG21	1:B:339:TRP:NE1	2.12	0.64
1:D:22:ILE:HD13	1:D:149:PRO:HB2	1.80	0.64
1:B:291:VAL:HG21	1:B:339:TRP:CD1	2.33	0.63
1:B:120:TRP:HE1	1:B:132:GLN:HG2	1.64	0.63
1:B:113:GLU:HG3	4:B:2029:HOH:O	1.98	0.63
1:C:427:CYS:O	1:C:466:ALA:HB2	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:427:CYS:N	4:C:2105:HOH:O	2.31	0.63
1:A:14:PHE:CD2	1:A:220:ASP:HB3	2.34	0.63
1:B:120:TRP:CE2	1:B:132:GLN:HG2	2.34	0.62
1:D:218:MET:HE3	1:D:224:LEU:HB3	1.82	0.62
1:C:329:GLN:HE21	1:D:128:ARG:HH11	1.48	0.62
1:D:137:TRP:CE3	1:D:159:GLN:HB2	2.34	0.62
1:A:232:LEU:HD11	1:A:258:SER:HB2	1.81	0.62
1:B:23:GLN:HG2	4:B:2008:HOH:O	2.00	0.62
1:C:428:GLY:N	4:C:2105:HOH:O	2.06	0.62
1:C:16:GLU:HG3	1:C:20:GLU:OE1	2.01	0.61
1:A:304:VAL:HG13	1:A:350:LEU:HD22	1.82	0.61
1:B:315:LEU:HD21	1:B:381:LEU:HD12	1.83	0.61
1:D:198:LEU:HD22	1:D:449:PHE:CE2	2.35	0.61
1:B:58:PHE:HE2	1:B:159:GLN:HG3	1.65	0.61
1:C:426:HIS:CG	4:C:2105:HOH:O	2.53	0.60
1:B:234:GLU:OE1	1:B:238:ARG:NH1	2.35	0.60
1:A:40:ARG:HD2	4:A:2007:HOH:O	2.00	0.60
1:A:168:ARG:HG3	1:A:231:TYR:OH	2.01	0.60
1:A:178:MET:HG3	1:A:259:TRP:CE2	2.37	0.60
1:A:315:LEU:HD21	1:A:381:LEU:HD12	1.81	0.60
1:C:460:ARG:HG3	1:C:460:ARG:HH11	1.67	0.60
1:B:35:GLY:HA2	1:B:147:THR:OG1	2.02	0.60
1:C:74:MET:HE2	1:C:107:ALA:HB2	1.84	0.59
1:C:120:TRP:CE2	1:C:132:GLN:HG2	2.37	0.59
1:C:198:LEU:HD22	1:C:449:PHE:CE2	2.37	0.59
1:D:370:GLU:OE1	1:D:372:LEU:HD21	2.02	0.59
1:A:9:PRO:HG2	1:A:12:GLN:HB2	1.84	0.59
1:A:309:GLN:HG2	1:A:312:ARG:NH2	2.18	0.59
1:B:94:VAL:HG13	1:B:106:LEU:HB2	1.82	0.59
1:D:72:ILE:HG12	1:D:115:VAL:CG2	2.33	0.59
1:D:9:PRO:HG2	1:D:12:GLN:CB	2.25	0.59
1:A:68:PHE:CD2	1:A:116:ILE:HG12	2.38	0.59
1:B:349:GLU:HG3	1:B:361:VAL:HB	1.85	0.59
1:A:137:TRP:CE3	1:A:159:GLN:HB2	2.38	0.58
1:C:51:GLN:HG2	1:C:137:TRP:CZ3	2.38	0.58
1:A:56:ARG:O	1:A:60:GLU:HG3	2.03	0.58
1:A:329:GLN:HB2	1:B:128:ARG:HA	1.85	0.58
1:B:12:GLN:CG	1:D:12:GLN:CG	2.81	0.58
1:C:58:PHE:HE2	1:C:159:GLN:HG3	1.69	0.58
1:C:76:PHE:HD2	1:C:114:THR:HB	1.68	0.58
1:A:53:VAL:HG12	1:A:57:MET:CE	2.33	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:VAL:CG1	1:A:57:MET:HE2	2.32	0.57
1:A:120:TRP:CE2	1:A:132:GLN:HG2	2.39	0.57
1:B:137:TRP:CZ3	1:B:159:GLN:HB2	2.39	0.57
1:D:309:GLN:HG2	1:D:312:ARG:NH2	2.20	0.57
1:D:435:LEU:HD12	1:D:460:ARG:NH2	2.18	0.57
1:A:214:ILE:HG13	1:A:214:ILE:O	2.05	0.57
1:B:178:MET:CE	1:B:181:ILE:HD12	2.35	0.57
1:C:72:ILE:HG12	1:C:115:VAL:CG2	2.35	0.57
1:B:26:GLU:HA	4:B:2009:HOH:O	2.05	0.56
1:C:336:PHE:HD1	1:C:347:ARG:HD2	1.70	0.56
1:D:118:TYR:O	1:D:121:SER:HB3	2.04	0.56
1:D:397:LEU:O	1:D:401:GLU:HG3	2.05	0.56
1:A:95:THR:OG1	1:A:96:HIS:HD2	1.88	0.56
1:B:137:TRP:CE3	1:B:159:GLN:HB2	2.40	0.56
1:B:156:PHE:HA	1:B:263:TRP:CZ2	2.40	0.56
1:D:205:PHE:HB2	1:D:230:HIS:CE1	2.40	0.56
1:A:74:MET:HE1	1:A:93:VAL:HG22	1.86	0.56
1:B:90:GLU:HB3	4:B:2031:HOH:O	2.06	0.56
1:C:214:ILE:HG13	1:C:214:ILE:O	2.05	0.56
1:C:426:HIS:HE1	1:C:433:GLU:HB2	1.70	0.56
1:D:248:ASP:O	1:D:250:GLN:HG3	2.05	0.56
1:D:296:TYR:HB2	1:D:300:SER:HB2	1.87	0.56
1:C:94:VAL:CG1	1:C:106:LEU:HB2	2.36	0.55
1:C:209:VAL:HG13	1:C:231:TYR:O	2.06	0.55
1:D:14:PHE:CD2	1:D:220:ASP:HB3	2.41	0.55
1:B:199:LYS:HE2	1:B:213:THR:HG21	1.89	0.55
1:A:309:GLN:HG2	1:A:312:ARG:HH22	1.71	0.55
1:B:18:TYR:O	1:B:21:VAL:HG12	2.07	0.55
1:A:233:GLY:O	1:A:255:HIS:HA	2.06	0.55
1:B:373:PRO:HG2	1:B:376:ALA:HB3	1.87	0.55
1:B:54:LEU:HD21	1:B:185:LEU:HD22	1.88	0.55
1:A:135:ASN:OD1	1:A:159:GLN:HG2	2.07	0.54
1:C:329:GLN:HE21	1:D:128:ARG:NH1	2.06	0.54
1:B:42:TYR:CE1	1:B:342:LYS:HB3	2.43	0.54
1:B:36:THR:HG21	4:B:2012:HOH:O	2.08	0.54
1:C:150:PHE:O	1:C:264:ARG:HA	2.08	0.54
1:D:53:VAL:HG12	1:D:57:MET:HE3	1.90	0.54
1:D:178:MET:HG3	1:D:259:TRP:CE2	2.42	0.54
1:B:22:ILE:HD13	1:B:149:PRO:HB2	1.89	0.54
1:B:12:GLN:CG	1:D:12:GLN:HG3	2.35	0.54
1:D:56:ARG:O	1:D:60:GLU:HG3	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:51:GLN:HA	1:C:137:TRP:CH2	2.43	0.54
1:A:72:ILE:HG12	1:A:115:VAL:CG1	2.37	0.53
1:B:315:LEU:CD2	1:B:381:LEU:HD12	2.38	0.53
1:A:388:PHE:CE2	1:A:392:LEU:HD11	2.43	0.53
1:C:297:LYS:H	1:C:300:SER:HB2	1.72	0.53
1:A:68:PHE:CE2	1:A:116:ILE:HG12	2.43	0.53
1:B:277:ARG:HG3	1:B:277:ARG:HH11	1.73	0.53
1:D:211:THR:HG23	1:D:230:HIS:ND1	2.24	0.53
1:C:137:TRP:CZ3	1:C:159:GLN:HB2	2.43	0.53
1:C:25:ALA:HB3	1:C:27:LEU:HD13	1.91	0.53
1:C:22:ILE:HD13	1:C:149:PRO:HB2	1.91	0.53
1:D:18:TYR:CZ	1:D:22:ILE:HD11	2.44	0.53
1:C:40:ARG:HD3	1:D:123:TRP:CZ2	2.44	0.53
1:C:137:TRP:CE3	1:C:159:GLN:HB2	2.45	0.52
1:D:97:ALA:HB3	1:D:102:LEU:HD21	1.90	0.52
1:B:15:SER:HB3	1:B:148:ARG:HH22	1.74	0.52
1:B:457:PHE:HB3	1:B:462:GLY:O	2.10	0.52
1:A:40:ARG:HD3	1:B:123:TRP:CZ2	2.44	0.52
1:A:473:PHE:O	1:A:474:ALA:HB2	2.09	0.52
1:B:12:GLN:CG	1:D:12:GLN:HG2	2.40	0.52
1:A:65:ASN:HA	1:A:135:ASN:O	2.10	0.52
1:A:137:TRP:CZ3	1:A:159:GLN:HB2	2.44	0.52
1:B:232:LEU:HD11	1:B:258:SER:HB2	1.91	0.52
1:D:199:LYS:HE2	1:D:213:THR:HG21	1.92	0.52
1:D:231:TYR:CZ	1:D:233:GLY:HA2	2.44	0.52
1:B:15:SER:HA	1:B:148:ARG:NH2	2.25	0.52
1:B:178:MET:HG3	1:B:259:TRP:CE2	2.45	0.52
1:B:57:MET:SD	1:B:181:ILE:HG23	2.50	0.52
1:B:331:THR:HG23	1:B:334:TYR:H	1.75	0.52
1:C:460:ARG:HD2	1:C:460:ARG:O	2.10	0.52
1:A:72:ILE:HG12	1:A:115:VAL:HG13	1.91	0.51
1:B:169:GLU:CD	1:B:169:GLU:H	2.18	0.51
1:D:111:THR:HG1	2:D:1478:PRO:N	2.08	0.51
1:A:132:GLN:O	1:A:163:THR:HA	2.11	0.51
1:B:26:GLU:O	1:B:40:ARG:HB2	2.10	0.51
1:C:258:SER:OG	2:C:1478:PRO:HB3	2.10	0.51
1:B:309:GLN:HG2	1:B:312:ARG:NH2	2.26	0.51
1:B:410:TYR:O	1:B:414:LYS:HG3	2.10	0.51
1:A:238:ARG:HB3	1:A:238:ARG:NH1	2.26	0.51
1:A:295:ILE:O	1:A:352:PRO:HD3	2.11	0.51
1:A:262:SER:C	1:A:264:ARG:H	2.19	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:53:VAL:HG12	1:C:57:MET:HE2	1.92	0.51
1:A:336:PHE:CE1	1:A:349:GLU:HG2	2.45	0.50
1:B:97:ALA:HB3	1:B:102:LEU:HD21	1.92	0.50
1:A:69:PRO:HD3	1:B:38:VAL:HG12	1.93	0.50
1:B:169:GLU:O	1:B:173:GLU:HG3	2.12	0.50
1:A:315:LEU:CD2	1:A:381:LEU:HD12	2.41	0.50
1:B:296:TYR:CD1	1:B:296:TYR:C	2.89	0.50
1:B:437:GLN:O	1:B:441:THR:HA	2.12	0.50
1:C:178:MET:HE2	1:C:259:TRP:CH2	2.47	0.50
1:B:234:GLU:O	1:B:238:ARG:HG3	2.12	0.50
1:D:329:GLN:H	1:D:329:GLN:NE2	2.09	0.50
1:D:72:ILE:HG12	1:D:115:VAL:HG21	1.93	0.50
1:D:249:LEU:N	1:D:249:LEU:HD22	2.27	0.50
1:D:137:TRP:CZ3	1:D:159:GLN:HB2	2.47	0.50
1:D:309:GLN:HG2	1:D:312:ARG:HH22	1.77	0.50
1:A:211:THR:HG23	1:A:230:HIS:CE1	2.47	0.50
1:A:437:GLN:O	1:A:441:THR:N	2.42	0.50
1:B:74:MET:CE	1:B:107:ALA:HB2	2.36	0.50
1:C:72:ILE:HG12	1:C:115:VAL:HG21	1.94	0.50
1:A:238:ARG:HA	1:A:253:TYR:CE2	2.47	0.50
1:C:15:SER:CA	1:C:148:ARG:NH2	2.73	0.49
1:A:205:PHE:HB2	1:A:230:HIS:CE1	2.47	0.49
1:C:357:GLY:O	1:C:359:GLN:HG3	2.12	0.49
1:D:58:PHE:HE2	1:D:159:GLN:HG3	1.77	0.49
1:D:158:TRP:CD1	1:D:158:TRP:C	2.90	0.49
1:B:150:PHE:O	1:B:264:ARG:HA	2.12	0.49
1:B:176:ARG:HD2	4:B:2053:HOH:O	2.13	0.49
1:B:455:GLU:HG3	4:B:2108:HOH:O	2.12	0.49
1:A:89:PRO:O	1:B:98:GLY:HA2	2.11	0.49
1:C:326:ASP:O	1:C:327:ARG:C	2.55	0.49
1:D:331:THR:HG23	1:D:334:TYR:HB2	1.94	0.49
1:D:370:GLU:HG3	1:D:372:LEU:HG	1.95	0.49
1:C:113:GLU:OE1	2:C:1478:PRO:N	2.45	0.49
1:A:192:ILE:HG12	1:A:272:THR:HG21	1.94	0.49
1:C:247:ARG:HG3	1:C:247:ARG:HH11	1.77	0.49
1:D:96:HIS:CE1	1:D:101:GLU:HG3	2.48	0.49
1:A:128:ARG:HH12	1:B:329:GLN:HB2	1.77	0.49
1:B:327:ARG:NH2	4:B:2084:HOH:O	2.45	0.49
1:A:133:LEU:HA	1:A:163:THR:HG22	1.94	0.49
1:B:440:THR:O	1:B:441:THR:OG1	2.27	0.49
1:C:202:LYS:HE3	4:C:2108:HOH:O	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:132:GLN:O	1:D:163:THR:HA	2.13	0.49
1:A:57:MET:SD	1:A:181:ILE:HG23	2.53	0.49
1:C:158:TRP:CD1	1:C:158:TRP:C	2.91	0.49
1:C:366:LEU:HD13	1:C:388:PHE:CZ	2.47	0.49
1:A:178:MET:HG3	1:A:259:TRP:CD2	2.48	0.48
1:D:295:ILE:O	1:D:352:PRO:HD3	2.13	0.48
1:B:198:LEU:CD2	1:B:449:PHE:HE2	2.24	0.48
1:D:33:VAL:O	1:D:36:THR:HG22	2.13	0.48
1:A:246:ASP:HB3	1:A:248:ASP:OD1	2.13	0.48
1:D:473:PHE:O	1:D:474:ALA:HB2	2.12	0.48
1:B:346:PHE:HB3	1:B:362:LEU:HD11	1.95	0.48
1:C:241:ASP:OD1	1:C:243:LYS:NZ	2.36	0.48
1:B:179:LEU:HD22	1:B:212:THR:HB	1.96	0.48
1:C:178:MET:HG3	1:C:259:TRP:CE2	2.48	0.48
1:A:169:GLU:CD	1:A:169:GLU:H	2.22	0.47
1:B:262:SER:C	1:B:264:ARG:H	2.22	0.47
1:D:133:LEU:HA	1:D:163:THR:HG22	1.96	0.47
1:A:388:PHE:HE2	1:A:392:LEU:HD11	1.79	0.47
1:B:12:GLN:HG2	1:D:12:GLN:HG2	1.92	0.47
1:C:426:HIS:CE1	1:C:433:GLU:HB2	2.49	0.47
1:B:15:SER:CA	1:B:148:ARG:NH2	2.77	0.47
1:B:152:ARG:HH11	1:B:152:ARG:HG3	1.79	0.47
1:B:214:ILE:HG13	1:B:214:ILE:O	2.14	0.47
1:C:135:ASN:OD1	1:C:159:GLN:HG2	2.15	0.47
1:B:399:PHE:O	1:B:403:HIS:HD2	1.97	0.47
1:D:120:TRP:CE2	1:D:132:GLN:HG2	2.47	0.47
1:D:435:LEU:HD12	1:D:439:GLU:OE1	2.14	0.47
1:A:158:TRP:CD1	1:A:158:TRP:C	2.93	0.47
1:A:221:GLY:O	1:A:420:GLY:HA2	2.15	0.47
1:B:336:PHE:HD1	1:B:347:ARG:HD2	1.80	0.47
1:C:426:HIS:CE1	4:C:2105:HOH:O	2.62	0.46
1:A:51:GLN:HA	1:A:137:TRP:CH2	2.50	0.46
1:A:15:SER:CA	1:A:148:ARG:HH22	2.27	0.46
1:B:224:LEU:HD12	1:B:264:ARG:HH21	1.80	0.46
1:D:234:GLU:CD	1:D:238:ARG:HH12	2.23	0.46
1:A:445:ARG:O	1:A:446:CYS:HB3	2.15	0.46
1:B:336:PHE:CD1	1:B:347:ARG:HD2	2.50	0.46
1:C:311:LEU:O	1:C:315:LEU:HG	2.14	0.46
1:D:296:TYR:HB3	1:D:304:VAL:HG21	1.97	0.46
1:A:381:LEU:O	1:A:382:PRO:C	2.56	0.46
1:B:349:GLU:CG	1:B:361:VAL:HB	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:390:GLU:HG3	4:D:2110:HOH:O	2.14	0.46
1:B:304:VAL:HG13	1:B:350:LEU:HD22	1.97	0.46
1:C:247:ARG:HG3	1:C:247:ARG:NH1	2.29	0.46
1:C:42:TYR:CE1	1:C:342:LYS:HB3	2.50	0.46
1:B:51:GLN:HA	1:B:137:TRP:CH2	2.51	0.46
1:B:473:PHE:O	1:B:474:ALA:HB2	2.16	0.46
1:C:74:MET:CE	1:C:107:ALA:HB2	2.45	0.46
1:C:234:GLU:CD	4:C:2070:HOH:O	2.59	0.46
1:C:296:TYR:CD1	1:C:296:TYR:C	2.94	0.46
1:A:247:ARG:HG3	1:A:247:ARG:HH11	1.81	0.46
1:B:378:PRO:HB2	4:B:2091:HOH:O	2.15	0.45
1:C:40:ARG:HD2	4:C:2015:HOH:O	2.15	0.45
1:A:128:ARG:NH1	1:B:329:GLN:HB2	2.30	0.45
1:C:249:LEU:N	1:C:249:LEU:CD1	2.79	0.45
1:C:14:PHE:CD2	1:C:220:ASP:HB3	2.51	0.45
1:B:377:LEU:CB	1:B:378:PRO:HD3	2.46	0.45
1:B:381:LEU:HB2	1:B:382:PRO:HD3	1.98	0.45
1:C:113:GLU:H	1:C:113:GLU:HG2	1.52	0.45
1:D:14:PHE:CG	1:D:220:ASP:HB3	2.51	0.45
1:A:93:VAL:O	1:B:94:VAL:HG23	2.17	0.45
1:A:110:PRO:HG2	1:A:111:THR:HG22	1.99	0.45
1:A:217:LEU:C	1:A:218:MET:HE2	2.42	0.45
1:A:304:VAL:HG13	1:A:350:LEU:CD2	2.46	0.45
1:D:135:ASN:OD1	1:D:159:GLN:HG2	2.16	0.45
1:D:179:LEU:HD22	1:D:212:THR:HB	1.99	0.45
1:C:194:VAL:HG12	1:C:216:ALA:HB2	1.98	0.45
1:B:202:LYS:HD3	1:B:433:GLU:OE2	2.17	0.45
1:C:74:MET:HE1	1:C:91:LEU:HD23	1.99	0.45
1:D:329:GLN:NE2	1:D:329:GLN:N	2.65	0.45
1:B:381:LEU:HA	1:B:381:LEU:HD23	1.75	0.45
1:C:205:PHE:HB2	1:C:230:HIS:CE1	2.52	0.45
1:A:42:TYR:CZ	1:A:342:LYS:HB3	2.53	0.44
1:A:198:LEU:CD2	1:A:449:PHE:HE2	2.22	0.44
1:B:15:SER:HA	1:B:148:ARG:HH21	1.80	0.44
1:B:391:GLU:OE2	1:B:391:GLU:HA	2.17	0.44
1:B:373:PRO:HG2	1:B:376:ALA:CB	2.46	0.44
1:D:148:ARG:HB3	1:D:152:ARG:HB2	1.99	0.44
1:D:331:THR:O	1:D:334:TYR:HB3	2.16	0.44
1:D:370:GLU:CD	1:D:372:LEU:HD21	2.42	0.44
1:C:447:VAL:HG22	1:C:471:VAL:HG22	2.00	0.44
1:D:331:THR:HG23	1:D:334:TYR:CB	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:135:ASN:OD1	1:B:159:GLN:HG2	2.17	0.44
1:B:466:ALA:C	1:B:468:GLY:H	2.25	0.44
1:D:16:GLU:HG3	1:D:20:GLU:OE1	2.17	0.44
1:D:262:SER:C	1:D:264:ARG:H	2.26	0.44
1:A:113:GLU:OE1	2:A:1478:PRO:N	2.50	0.44
1:B:295:ILE:N	1:B:295:ILE:HD12	2.32	0.44
1:A:139:ASN:HB3	1:A:157:LEU:HD23	2.00	0.44
1:B:9:PRO:HB3	1:D:11:SER:HB2	2.00	0.44
1:B:175:VAL:HG13	1:B:229:SER:OG	2.17	0.44
1:A:97:ALA:HB1	1:B:143:TRP:CD2	2.52	0.44
1:A:305:LEU:HD21	1:A:326:ASP:OD1	2.17	0.44
1:A:447:VAL:HG11	1:A:467:TYR:HB3	2.00	0.44
1:D:109:ARG:HG3	1:D:109:ARG:O	2.18	0.43
1:D:381:LEU:O	1:D:382:PRO:C	2.60	0.43
1:A:94:VAL:HA	1:B:94:VAL:HA	2.00	0.43
1:C:22:ILE:CD1	1:C:149:PRO:HB2	2.48	0.43
1:C:459:VAL:HG22	1:C:459:VAL:O	2.17	0.43
1:D:232:LEU:HD11	1:D:258:SER:HB2	2.00	0.43
1:D:238:ARG:HA	1:D:253:TYR:CE2	2.53	0.43
1:A:231:TYR:CE1	1:A:233:GLY:HA2	2.53	0.43
1:A:366:LEU:HD12	1:A:366:LEU:HA	1.78	0.43
1:A:399:PHE:O	1:A:403:HIS:HD2	2.00	0.43
1:C:36:THR:HG21	4:C:2011:HOH:O	2.18	0.43
1:D:366:LEU:HD22	1:D:388:PHE:CE1	2.53	0.43
1:C:277:ARG:HG2	1:C:277:ARG:HH11	1.83	0.43
1:C:318:GLN:HG3	1:C:378:PRO:O	2.18	0.43
1:B:283:PRO:O	1:B:321:ARG:NH2	2.51	0.43
1:B:311:LEU:O	1:B:312:ARG:C	2.61	0.43
1:C:45:ALA:HB1	1:C:287:PRO:HB3	1.98	0.43
1:B:395:ARG:O	1:B:398:ALA:HB3	2.19	0.43
1:C:65:ASN:HA	1:C:135:ASN:O	2.19	0.43
1:D:209:VAL:CG2	1:D:231:TYR:HD2	2.32	0.43
1:D:214:ILE:O	1:D:214:ILE:HG13	2.18	0.43
1:D:373:PRO:HG2	1:D:376:ALA:HB3	1.99	0.43
1:A:451:ALA:HB3	1:A:469:LYS:HD3	2.01	0.43
1:A:358:GLY:O	1:A:373:PRO:HA	2.19	0.43
1:C:266:ILE:O	1:C:270:ILE:HG13	2.18	0.43
1:D:198:LEU:CD2	1:D:449:PHE:HE2	2.29	0.43
1:C:74:MET:CE	1:C:93:VAL:HG22	2.49	0.42
1:D:18:TYR:O	1:D:21:VAL:HG12	2.19	0.42
1:A:201:GLU:HG3	4:A:2087:HOH:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:291:VAL:HG21	1:A:339:TRP:CD1	2.54	0.42
1:B:54:LEU:HD23	1:B:57:MET:HE3	2.01	0.42
1:D:73:PRO:O	1:D:76:PHE:HB2	2.19	0.42
1:D:291:VAL:HG13	1:D:347:ARG:HG2	2.00	0.42
1:D:410:TYR:CE2	1:D:460:ARG:HD3	2.53	0.42
1:D:460:ARG:HG3	1:D:460:ARG:NH1	2.28	0.42
1:A:178:MET:CE	1:A:181:ILE:HD12	2.50	0.42
1:A:425:PHE:CZ	1:A:454:GLU:HB3	2.54	0.42
1:B:113:GLU:H	1:B:113:GLU:HG2	1.36	0.42
1:B:283:PRO:HB3	1:B:385:LEU:HD12	2.02	0.42
1:B:339:TRP:HA	1:B:339:TRP:CE3	2.54	0.42
1:A:15:SER:O	1:A:19:LEU:HD22	2.20	0.42
1:A:329:GLN:CB	1:B:128:ARG:HA	2.50	0.42
1:B:366:LEU:HD13	1:B:388:PHE:CZ	2.55	0.42
1:A:247:ARG:HG3	1:A:247:ARG:NH1	2.34	0.42
1:C:373:PRO:HG2	1:C:376:ALA:HB3	2.00	0.42
1:D:150:PHE:O	1:D:264:ARG:HA	2.18	0.42
1:C:410:TYR:O	1:C:414:LYS:HG3	2.20	0.42
1:D:437:GLN:O	1:D:441:THR:HA	2.19	0.42
1:A:116:ILE:HD12	1:A:162:HIS:CD2	2.55	0.42
1:B:15:SER:HB3	1:B:148:ARG:NH2	2.34	0.42
1:B:65:ASN:HA	1:B:135:ASN:O	2.20	0.42
1:B:94:VAL:CG1	1:B:106:LEU:HB2	2.47	0.42
1:B:100:GLU:H	1:B:100:GLU:HG3	1.66	0.42
1:B:327:ARG:HB3	1:B:329:GLN:OE1	2.19	0.42
1:C:120:TRP:HE1	1:C:132:GLN:CG	2.25	0.42
1:D:445:ARG:O	1:D:446:CYS:HB3	2.19	0.42
1:C:232:LEU:HD11	1:C:258:SER:HB2	2.02	0.42
1:D:360:ALA:HB2	1:D:374:LEU:HD21	2.01	0.42
1:C:54:LEU:HD23	1:C:57:MET:HE3	2.02	0.42
1:D:201:GLU:HB3	1:D:430:LYS:HE2	2.00	0.42
1:A:353:LYS:HA	1:A:353:LYS:HD3	1.76	0.41
1:B:300:SER:O	1:B:304:VAL:HG23	2.20	0.41
1:B:341:LEU:HD23	1:B:341:LEU:HA	1.86	0.41
1:C:437:GLN:O	1:C:441:THR:HA	2.20	0.41
1:A:94:VAL:HG13	1:A:106:LEU:HB2	2.02	0.41
1:B:231:TYR:CZ	1:B:233:GLY:HA2	2.54	0.41
1:B:360:ALA:O	1:B:371:THR:HA	2.20	0.41
1:C:40:ARG:HB3	1:C:41:PRO:HD2	2.01	0.41
1:C:76:PHE:CD2	1:C:114:THR:HB	2.52	0.41
1:C:156:PHE:HA	1:C:263:TRP:CZ2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:460:ARG:HG3	1:C:460:ARG:NH1	2.33	0.41
1:D:435:LEU:HD12	1:D:460:ARG:HH21	1.83	0.41
1:A:291:VAL:CG1	1:A:347:ARG:HG3	2.50	0.41
1:B:261:LEU:HD23	1:B:261:LEU:C	2.45	0.41
1:B:296:TYR:HE2	1:B:305:LEU:HD21	1.84	0.41
1:B:184:ARG:HG3	1:B:188:GLU:OE2	2.21	0.41
1:B:288:ILE:O	1:B:344:VAL:HG13	2.20	0.41
1:D:40:ARG:HB3	1:D:41:PRO:HD2	2.03	0.41
1:D:277:ARG:NH1	1:D:340:GLU:OE2	2.53	0.41
1:D:291:VAL:HG13	1:D:291:VAL:O	2.20	0.41
1:B:11:SER:HB2	1:D:9:PRO:HB3	2.02	0.41
1:B:14:PHE:O	1:B:17:TRP:HB3	2.21	0.41
1:C:238:ARG:CD	4:C:2070:HOH:O	2.57	0.41
1:C:277:ARG:HG2	1:C:277:ARG:NH1	2.36	0.41
1:C:473:PHE:O	1:C:474:ALA:HB2	2.20	0.41
1:D:18:TYR:CE1	1:D:22:ILE:HD11	2.55	0.41
1:A:120:TRP:HB3	1:A:124:ILE:HD12	2.03	0.41
1:A:282:PRO:HA	1:A:283:PRO:HD3	1.87	0.41
1:B:205:PHE:HB2	1:B:230:HIS:CE1	2.55	0.41
1:D:373:PRO:HG2	1:D:376:ALA:CB	2.51	0.41
1:A:51:GLN:O	1:A:55:ASP:HB2	2.21	0.41
1:A:94:VAL:HG23	1:B:93:VAL:O	2.21	0.41
1:A:115:VAL:HG23	1:A:116:ILE:N	2.35	0.41
1:B:88:SER:HA	1:B:89:PRO:HD3	1.87	0.41
1:D:443:THR:HA	4:D:2118:HOH:O	2.21	0.41
1:D:460:ARG:NH1	1:D:460:ARG:CG	2.83	0.41
1:A:45:ALA:HB1	1:A:287:PRO:HB3	2.03	0.40
1:A:437:GLN:O	1:A:441:THR:HA	2.22	0.40
1:C:179:LEU:HD22	1:C:212:THR:HB	2.03	0.40
1:C:264:ARG:CG	1:C:264:ARG:NH1	2.61	0.40
1:D:202:LYS:HD3	1:D:433:GLU:OE2	2.22	0.40
1:A:18:TYR:CZ	1:A:22:ILE:HD11	2.57	0.40
1:A:54:LEU:HD23	1:A:57:MET:HE3	2.04	0.40
1:A:61:THR:CG2	1:A:177:ARG:HD3	2.51	0.40
1:B:61:THR:CG2	1:B:177:ARG:HD3	2.51	0.40
1:B:460:ARG:HD3	1:B:460:ARG:C	2.45	0.40
1:D:336:PHE:CE1	1:D:349:GLU:HG2	2.56	0.40
1:B:178:MET:HE1	1:B:181:ILE:HD12	2.03	0.40
1:B:297:LYS:O	1:B:298:ASP:C	2.63	0.40
1:C:410:TYR:OH	1:C:440:THR:HG23	2.20	0.40
1:B:14:PHE:CG	1:B:220:ASP:HB3	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:275:ASP:OD2	1:A:275:ASP:C	2.64	0.40
1:C:25:ALA:CB	1:C:27:LEU:HD13	2.52	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	460/477 (96%)	430 (94%)	29 (6%)	1 (0%)	43	72
1	B	460/477 (96%)	427 (93%)	30 (6%)	3 (1%)	18	48
1	C	460/477 (96%)	437 (95%)	22 (5%)	1 (0%)	43	72
1	D	460/477 (96%)	440 (96%)	20 (4%)	0	100	100
All	All	1840/1908 (96%)	1734 (94%)	101 (6%)	5 (0%)	36	65

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	467	TYR
1	C	462	GLY
1	A	474	ALA
1	B	352	PRO
1	B	280	VAL

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	385/395 (98%)	362 (94%)	23 (6%)	17	47
1	B	385/395 (98%)	361 (94%)	24 (6%)	16	46
1	C	385/395 (98%)	361 (94%)	24 (6%)	16	46
1	D	385/395 (98%)	364 (94%)	21 (6%)	19	50
All	All	1540/1580 (98%)	1448 (94%)	92 (6%)	17	47

All (92) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	LEU
1	A	36	THR
1	A	38	VAL
1	A	40	ARG
1	A	55	ASP
1	A	91	LEU
1	A	94	VAL
1	A	100	GLU
1	A	147	THR
1	A	159	GLN
1	A	198	LEU
1	A	228	THR
1	A	259	TRP
1	A	299	GLU
1	A	329	GLN
1	A	353	LYS
1	A	366	LEU
1	A	377	LEU
1	A	385	LEU
1	A	418	GLN
1	A	423	LEU
1	A	435	LEU
1	A	455	GLU
1	B	12	GLN
1	B	38	VAL
1	B	41	PRO
1	B	55	ASP
1	B	75	SER
1	B	90	GLU
1	B	91	LEU
1	B	94	VAL

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Mol	Chain	Res	Type
1	B	100	GLU
1	B	113	GLU
1	B	153	THR
1	B	198	LEU
1	B	209	VAL
1	B	243	LYS
1	B	259	TRP
1	B	277	ARG
1	B	329	GLN
1	B	331	THR
1	B	341	LEU
1	B	366	LEU
1	B	377	LEU
1	B	385	LEU
1	B	435	LEU
1	B	455	GLU
1	C	38	VAL
1	C	40	ARG
1	C	55	ASP
1	C	75	SER
1	C	91	LEU
1	C	94	VAL
1	C	113	GLU
1	C	147	THR
1	C	159	GLN
1	C	198	LEU
1	C	209	VAL
1	C	234	GLU
1	C	249	LEU
1	C	250	GLN
1	C	259	TRP
1	C	264	ARG
1	C	329	GLN
1	C	353	LYS
1	C	366	LEU
1	C	377	LEU
1	C	385	LEU
1	C	418	GLN
1	C	423	LEU
1	C	455	GLU
1	D	12	GLN
1	D	27	LEU

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Mol	Chain	Res	Type
1	D	36	THR
1	D	38	VAL
1	D	55	ASP
1	D	94	VAL
1	D	149	PRO
1	D	198	LEU
1	D	209	VAL
1	D	259	TRP
1	D	329	GLN
1	D	331	THR
1	D	341	LEU
1	D	353	LYS
1	D	366	LEU
1	D	377	LEU
1	D	385	LEU
1	D	390	GLU
1	D	418	GLN
1	D	435	LEU
1	D	455	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	51	GLN
1	A	96	HIS
1	A	132	GLN
1	A	245	GLN
1	A	273	HIS
1	A	309	GLN
1	A	403	HIS
1	B	51	GLN
1	B	64	GLN
1	B	309	GLN
1	B	330	HIS
1	C	12	GLN
1	C	64	GLN
1	C	132	GLN
1	C	309	GLN
1	C	329	GLN
1	C	337	HIS
1	C	403	HIS

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Mol	Chain	Res	Type
1	D	51	GLN
1	D	96	HIS
1	D	132	GLN
1	D	225	GLN
1	D	250	GLN
1	D	309	GLN
1	D	329	GLN
1	D	337	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 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	$\# Z > 2$
2	PRO	A	1478	-	8,8,8	0.79	0	10,10,10	1.00	0
2	PRO	C	1478	-	8,8,8	0.86	0	10,10,10	0.90	0
2	PRO	B	1478	-	8,8,8	0.82	0	10,10,10	0.95	0
2	PRO	D	1478	-	8,8,8	0.79	0	10,10,10	0.87	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
2	PRO	A	1478	-	-	0/4/11/11	0/1/1/1
2	PRO	C	1478	-	-	0/4/11/11	0/1/1/1
2	PRO	B	1478	-	-	0/4/11/11	0/1/1/1
2	PRO	D	1478	-	-	0/4/11/11	0/1/1/1

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.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1478	PRO	1	0
2	C	1478	PRO	2	0
2	D	1478	PRO	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	464/477 (97%)	-0.26	2 (0%) 88 85	15, 39, 65, 83	8 (1%)
1	B	464/477 (97%)	-0.30	6 (1%) 75 67	13, 31, 62, 93	7 (1%)
1	C	464/477 (97%)	-0.53	2 (0%) 88 85	10, 27, 55, 71	6 (1%)
1	D	464/477 (97%)	-0.59	3 (0%) 85 81	11, 27, 52, 83	7 (1%)
All	All	1856/1908 (97%)	-0.42	13 (0%) 84 79	10, 31, 61, 93	28 (1%)

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	326	ASP	3.8
1	B	300	SER	3.3
1	B	296	TYR	2.9
1	A	250	GLN	2.7
1	B	299	GLU	2.4
1	D	87	PHE	2.3
1	B	328	ASP	2.2
1	D	450	GLU	2.2
1	D	145	MET	2.1
1	B	450	GLU	2.1
1	A	347	ARG	2.1
1	C	464	PRO	2.0
1	C	427	CYS	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	ZN	C	490	1/1	0.94	0.06	57,57,57,57	0
3	ZN	D	490	1/1	0.97	0.04	39,39,39,39	0
2	PRO	C	1478	8/8	0.98	0.06	18,19,20,21	0
2	PRO	D	1478	8/8	0.98	0.05	19,22,23,25	0
3	ZN	B	490	1/1	0.98	0.04	51,51,51,51	0
2	PRO	A	1478	8/8	0.98	0.07	27,29,30,32	0
2	PRO	B	1478	8/8	0.98	0.07	12,16,19,20	0
3	ZN	A	490	1/1	0.99	0.04	62,62,62,62	0

6.5 Other polymers [i](#)

There are no such residues in this entry.