



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 16, 2016 – 08:06 PM EST

PDB ID : 5K8N
Title : 5NAA-bound 5-nitroanthranilate aminohydrolase
Authors : Kalyoncu, S.
Deposited on : 2016-05-30
Resolution : 3.23 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20028320
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20028320

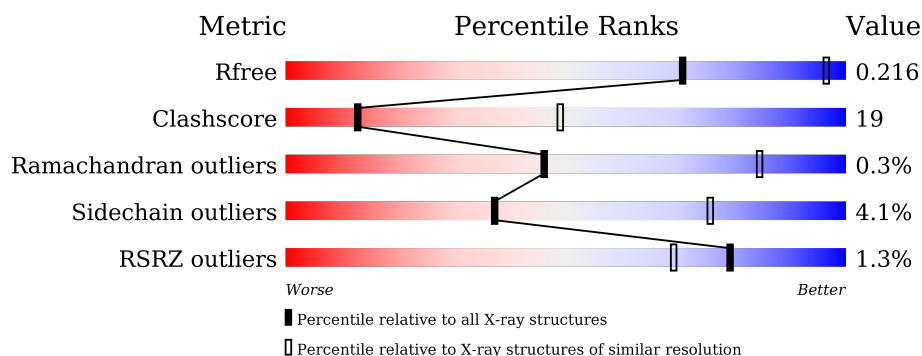
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.23 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}	91344	1095 (3.26-3.18)
Clashscore	102246	1046 (3.24-3.20)
Ramachandran outliers	100387	1026 (3.24-3.20)
Sidechain outliers	100360	1025 (3.24-3.20)
RSRZ outliers	91569	1100 (3.26-3.18)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	425	<div> <div></div> <div>60%36%..</div> </div>
1	B	425	<div> <div>2%</div> <div>58%40%..</div> </div>
1	C	425	<div> <div>5%</div> <div>62%35%..</div> </div>
1	D	425	<div> <div></div> <div>66%31%..</div> </div>
1	E	425	<div> <div></div> <div>64%33%..</div> </div>
1	F	425	<div> <div>%</div> <div>63%34%..</div> </div>

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Mol	Chain	Length	Quality of chain
1	G	425	<div> <div>%</div> <div> </div> <div>60%36%</div> <div>..</div> </div>
1	H	425	<div> <div>%</div> <div> </div> <div>63%34%</div> <div>..</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	6R6	A	501	-	-	-	X
2	6R6	D	501	-	-	-	X
2	6R6	F	501	-	-	-	X
2	6R6	G	501	-	-	-	X

2 Entry composition

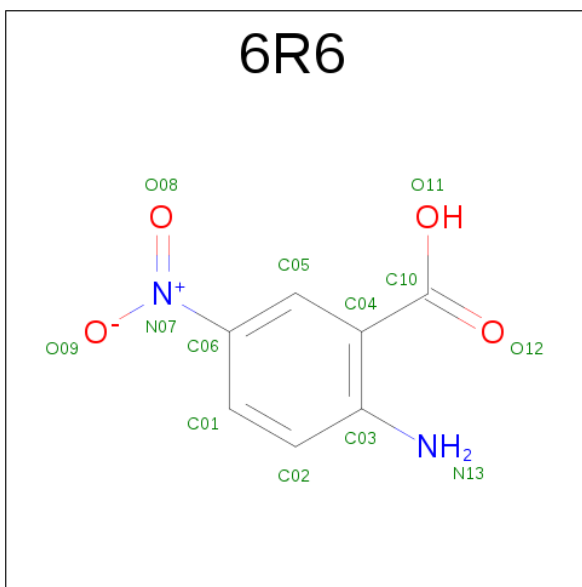
There are 2 unique types of molecules in this entry. The entry contains 25920 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 5-nitroanthranilic acid aminohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	420	Total	C	N	O	S	0	0	0
			3227	2042	560	606	19			
1	B	420	Total	C	N	O	S	0	0	0
			3227	2042	560	606	19			
1	C	420	Total	C	N	O	S	0	0	0
			3227	2042	560	606	19			
1	D	420	Total	C	N	O	S	0	0	0
			3227	2042	560	606	19			
1	E	420	Total	C	N	O	S	0	0	0
			3227	2042	560	606	19			
1	F	420	Total	C	N	O	S	0	0	0
			3227	2042	560	606	19			
1	G	420	Total	C	N	O	S	0	0	0
			3227	2042	560	606	19			
1	H	420	Total	C	N	O	S	0	0	0
			3227	2042	560	606	19			

- Molecule 2 is 5-nitroanthranilic acid (three-letter code: 6R6) (formula: C₇H₆N₂O₄).

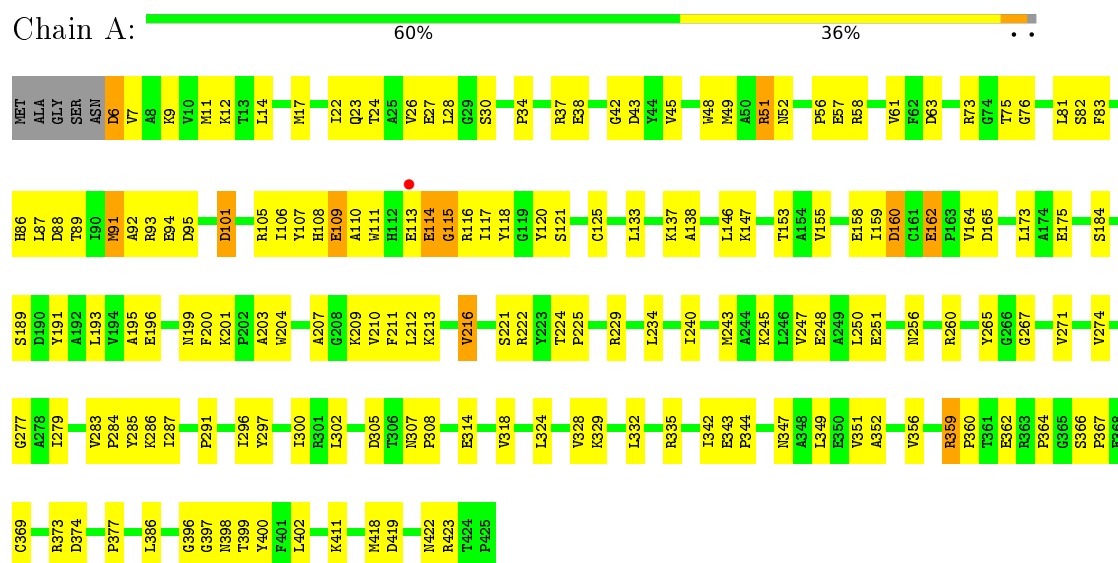


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			13	7	2	4		
2	B	1	Total	C	N	O	0	0
			13	7	2	4		
2	C	1	Total	C	N	O	0	0
			13	7	2	4		
2	D	1	Total	C	N	O	0	0
			13	7	2	4		
2	E	1	Total	C	N	O	0	0
			13	7	2	4		
2	F	1	Total	C	N	O	0	0
			13	7	2	4		
2	G	1	Total	C	N	O	0	0
			13	7	2	4		
2	H	1	Total	C	N	O	0	0
			13	7	2	4		

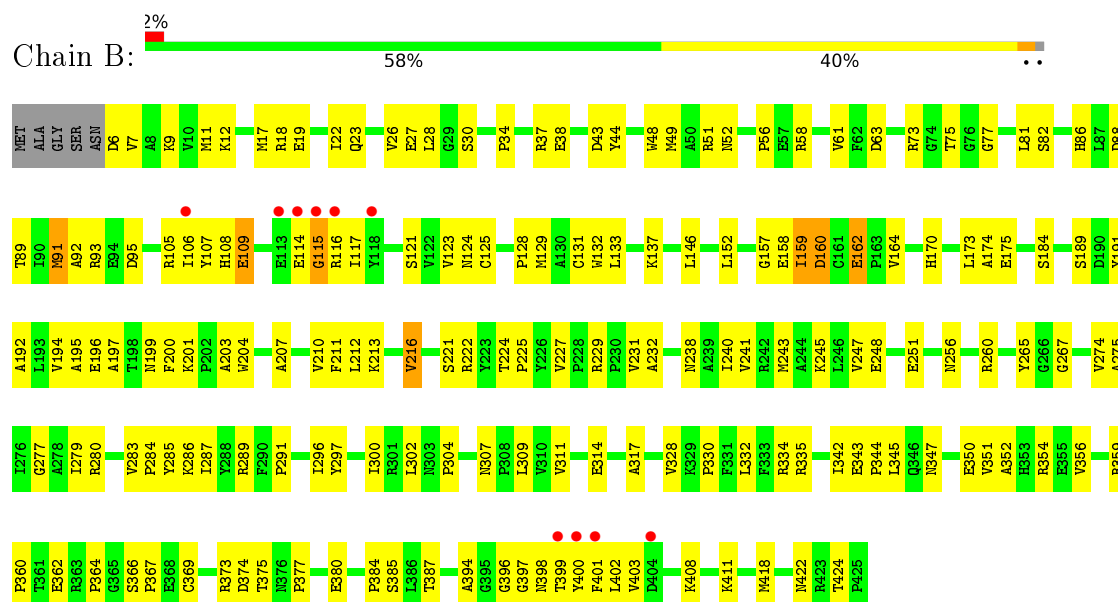
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

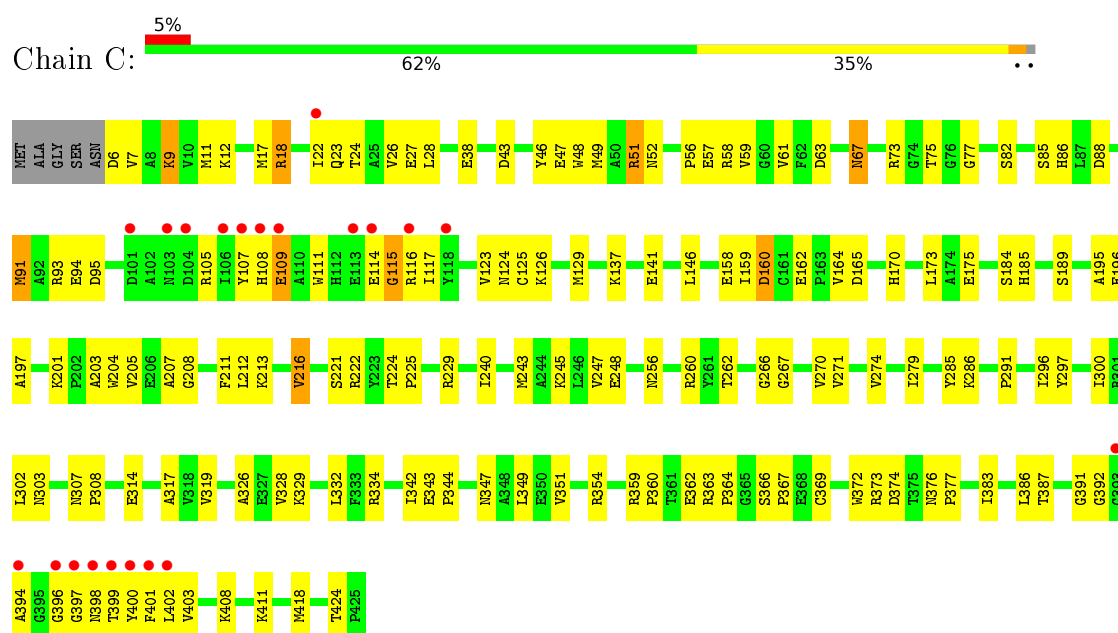
• Molecule 1: 5-nitroanthranilic acid aminohydrolase



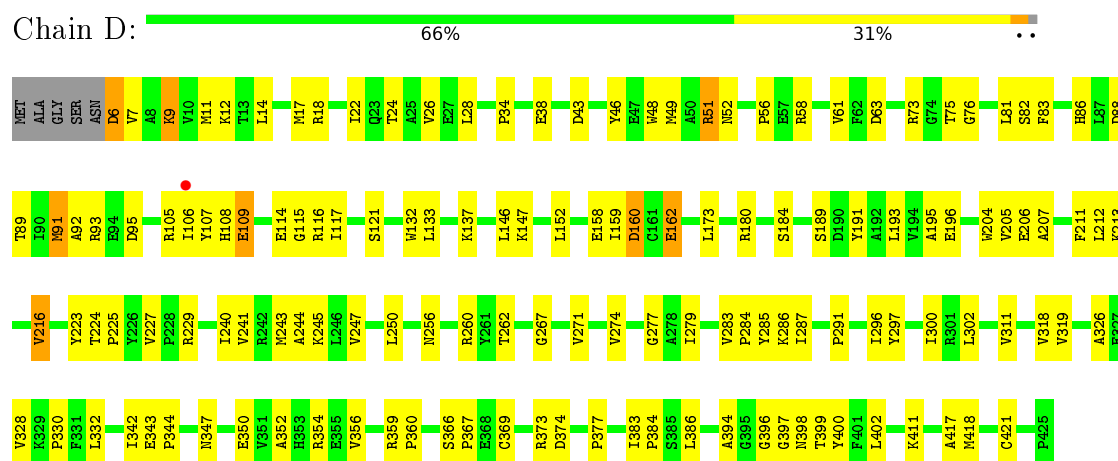
• Molecule 1: 5-nitroanthranilic acid aminohydrolase



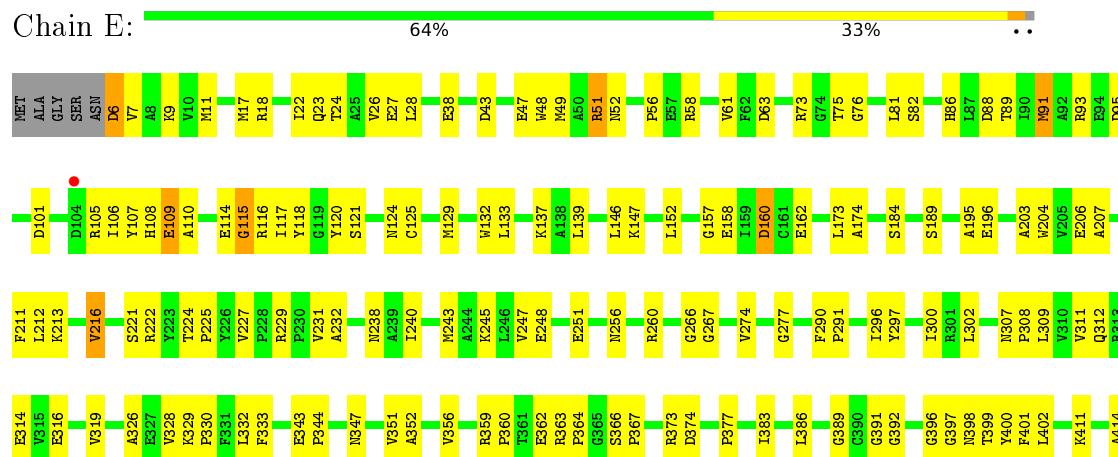
• Molecule 1: 5-nitroanthranilic acid aminohydrolase



- Molecule 1: 5-nitroanthranilic acid aminohydrolase

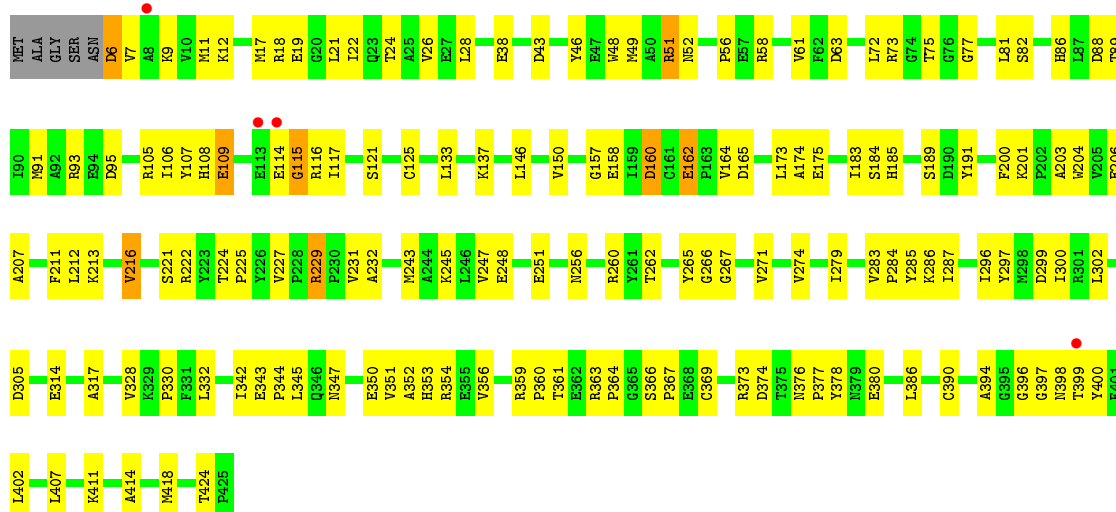


- Molecule 1: 5-nitroanthranilic acid aminohydrolase

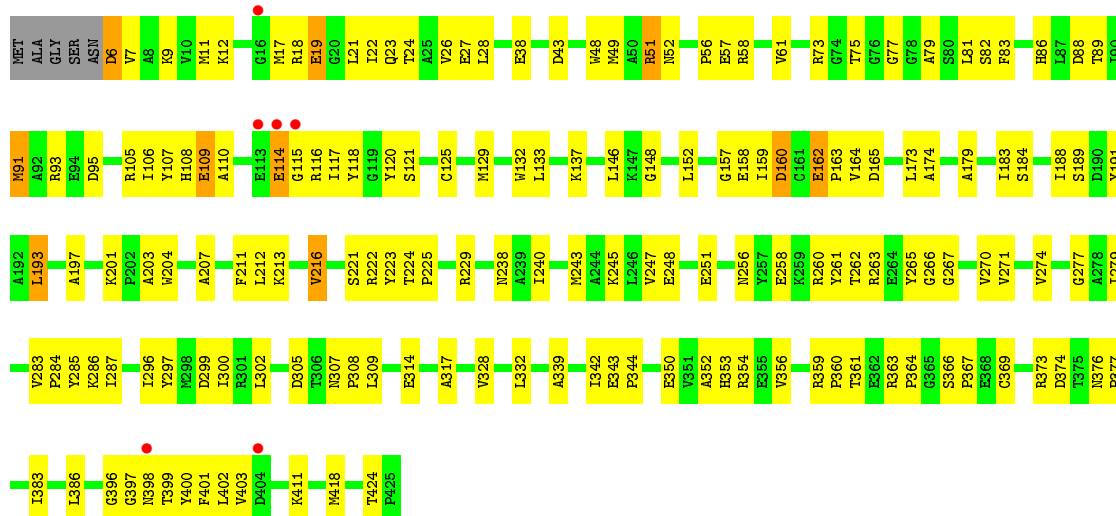




• Molecule 1: 5-nitroanthranilic acid aminohydrolase



• Molecule 1: 5-nitroanthranilic acid aminohydrolase



• Molecule 1: 5-nitroanthranilic acid aminohydrolase



R93	D190	I296	K411
E94	Y191	I297	
D95		I300	A417
	A195	I301	M418
D101	E196	L302	D419
			I420
R105	F200	D305	C421
I106		T306	M422
Y107	A207	N307	R423
H108		P308	T424
E109	F211	L309	P425
A110	L212		
	K213	E314	
E114			
G115	V216	V328	
R116			
I117	S221		
Y118	R222	L332	
G119	Y223		
Y120	T224	I342	
S121	P225	E343	
	Y226	P344	
N124	V227		
C125	P228	L349	
	R229		
L133		R359	
I134	I240	P360	
A135			
A136	M243	P364	
K137	A244	G365	
A138	K245	S366	
L139	L246	P367	
	V247	E368	
L146	E248	C369	
K147		S370	
	E252	N371	
V155	W253	G372	
G156		R373	
G157	N256	D374	
E158			
I159	R260	P377	
D160			
C161	R263	I383	
E162	E264	P384	
P163	Y265	S385	
V164		L386	
D165	V271	T387	
		Y388	
	V274	G389	
L173		C390	
A174	I279	G391	
E175			
		G396	
A179	Y285	G397	
R180	K286	N398	
	I287	T399	
I183	Y288	Y400	
S184	E289	F401	
H185	F290	L402	
	P291		
S189			

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	186.46Å 248.88Å 249.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.28 – 3.23 47.83 – 3.22	Depositor EDS
% Data completeness (in resolution range)	93.8 (46.28-3.23) 93.9 (47.83-3.22)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.39 (at 3.25Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.177 , 0.217 0.173 , 0.216	Depositor DCC
R_{free} test set	2001 reflections (2.29%)	DCC
Wilson B-factor (Å ²)	42.3	Xtriage
Anisotropy	0.750	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 32.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	25920	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.65% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 6R6

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	0.62	0/3301	0.64	0/4480
1	B	0.62	0/3301	0.66	0/4480
1	C	0.64	0/3301	0.68	0/4480
1	D	0.63	0/3301	0.65	0/4480
1	E	0.63	0/3301	0.64	0/4480
1	F	0.58	0/3301	0.64	0/4480
1	G	0.59	0/3301	0.64	0/4480
1	H	0.63	0/3301	0.66	0/4480
All	All	0.62	0/26408	0.65	0/35840

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	3227	0	3159	143	0
1	B	3227	0	3159	136	0
1	C	3227	0	3159	126	0
1	D	3227	0	3159	122	0
1	E	3227	0	3159	120	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	3227	0	3159	127	0
1	G	3227	0	3159	134	0
1	H	3227	0	3159	133	0
2	A	13	0	0	2	0
2	B	13	0	0	1	0
2	C	13	0	0	1	0
2	D	13	0	0	1	0
2	E	13	0	0	2	0
2	F	13	0	0	1	0
2	G	13	0	0	1	0
2	H	13	0	0	1	0
All	All	25920	0	25272	984	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 19.

All (984) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:61:VAL:HG13	1:H:173:LEU:HD11	1.36	1.08
1:D:173:LEU:HD11	1:F:61:VAL:HG13	1.37	1.02
1:A:49:MET:HB3	1:A:56:PRO:HG3	1.42	1.00
1:B:22:ILE:HG23	1:B:117:ILE:HD11	1.39	0.99
1:B:173:LEU:HD11	1:H:61:VAL:HG13	1.47	0.96
1:D:22:ILE:HG23	1:D:117:ILE:HD11	1.50	0.93
1:F:342:ILE:HD11	1:F:386:LEU:HD23	1.51	0.93
1:C:114:GLU:OE1	1:C:116:ARG:HG2	1.70	0.92
1:C:22:ILE:HG23	1:C:117:ILE:HD11	1.49	0.91
1:A:61:VAL:HG13	1:C:173:LEU:CD1	2.00	0.91
1:G:49:MET:HB3	1:G:56:PRO:HG3	1.51	0.91
1:F:203:ALA:HB2	1:F:364:PRO:HG3	1.52	0.91
1:G:22:ILE:HG23	1:G:117:ILE:HD11	1.54	0.88
1:D:212:LEU:HD12	1:D:300:ILE:CD1	2.03	0.88
1:B:173:LEU:CD1	1:H:61:VAL:HG13	2.04	0.88
1:H:49:MET:HB3	1:H:56:PRO:HG3	1.54	0.87
1:B:203:ALA:HB2	1:B:364:PRO:HG3	1.56	0.86
1:C:82:SER:HB2	1:C:189:SER:OG	1.76	0.86
1:E:243:MET:HE3	1:E:296:ILE:HG23	1.56	0.86
1:F:373:ARG:NH1	2:F:501:6R6:O11	2.08	0.86
1:A:6:ASP:HA	1:A:9:LYS:HE2	1.59	0.85
1:A:61:VAL:HG13	1:C:173:LEU:HD11	1.59	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:22:ILE:HG23	1:B:117:ILE:CD1	2.08	0.83
1:D:61:VAL:HG13	1:F:173:LEU:CD1	2.09	0.83
1:H:307:ASN:ND2	1:H:308:PRO:HD2	1.92	0.83
1:D:116:ARG:HD3	1:D:400:TYR:CD2	2.14	0.83
1:D:52:ASN:OD1	1:D:137:LYS:NZ	2.10	0.83
1:C:49:MET:HB3	1:C:56:PRO:HG3	1.59	0.82
1:B:82:SER:HB2	1:B:189:SER:OG	1.77	0.82
1:C:7:VAL:O	1:C:11:MET:HG2	1.79	0.81
1:A:212:LEU:HD12	1:A:300:ILE:HD12	1.62	0.81
1:G:243:MET:HE3	1:G:296:ILE:HG23	1.62	0.81
1:G:22:ILE:HG23	1:G:117:ILE:CD1	2.10	0.81
1:H:82:SER:HB2	1:H:189:SER:OG	1.81	0.81
1:C:212:LEU:HD12	1:C:300:ILE:HD12	1.62	0.81
1:D:114:GLU:O	1:D:116:ARG:N	2.13	0.80
1:B:114:GLU:OE1	1:B:116:ARG:HG2	1.81	0.80
1:A:38:GLU:HG2	1:A:88:ASP:HB3	1.61	0.80
1:A:173:LEU:HD11	1:C:61:VAL:HG13	1.63	0.80
1:A:7:VAL:HG22	1:A:418:MET:HE3	1.64	0.80
1:B:229:ARG:NH1	1:D:274:VAL:O	2.14	0.80
1:G:48:TRP:O	1:G:52:ASN:ND2	2.14	0.80
1:E:173:LEU:HD11	1:G:61:VAL:HG13	1.61	0.80
1:C:43:ASP:OD1	1:C:58:ARG:NH1	2.14	0.80
1:E:61:VAL:HG13	1:G:173:LEU:HD11	1.63	0.79
1:E:114:GLU:O	1:E:116:ARG:N	2.16	0.79
1:E:212:LEU:HD12	1:E:300:ILE:CD1	2.13	0.79
1:A:82:SER:HB2	1:A:189:SER:OG	1.82	0.79
1:F:22:ILE:HG23	1:F:117:ILE:HD11	1.64	0.79
1:H:7:VAL:HG22	1:H:418:MET:HE3	1.64	0.79
1:F:49:MET:HB3	1:F:56:PRO:HG3	1.65	0.79
1:H:114:GLU:O	1:H:116:ARG:N	2.15	0.79
1:D:82:SER:HB2	1:D:189:SER:OG	1.83	0.78
1:B:48:TRP:O	1:B:52:ASN:ND2	2.17	0.78
1:F:243:MET:HE3	1:F:296:ILE:HG23	1.64	0.78
1:C:398:ASN:OD1	1:C:399:THR:N	2.16	0.78
1:B:212:LEU:HD12	1:B:300:ILE:CD1	2.14	0.77
1:F:114:GLU:O	1:F:116:ARG:N	2.17	0.77
1:G:114:GLU:OE1	1:G:116:ARG:HG2	1.82	0.77
1:E:52:ASN:OD1	1:E:137:LYS:NZ	2.17	0.77
1:E:116:ARG:HD3	1:E:400:TYR:CD2	2.19	0.77
1:B:52:ASN:OD1	1:B:137:LYS:NZ	2.18	0.76
1:F:43:ASP:OD1	1:F:58:ARG:NH1	2.17	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:373:ARG:NH1	2:H:501:6R6:O12	2.18	0.76
1:E:61:VAL:HG13	1:G:173:LEU:CD1	2.15	0.76
1:A:48:TRP:O	1:A:52:ASN:ND2	2.18	0.76
1:B:7:VAL:O	1:B:11:MET:HG2	1.85	0.76
1:H:212:LEU:HD12	1:H:300:ILE:CD1	2.15	0.76
1:A:52:ASN:OD1	1:A:137:LYS:NZ	2.17	0.76
1:B:49:MET:HB3	1:B:56:PRO:HG3	1.67	0.76
1:F:82:SER:HB2	1:F:189:SER:OG	1.85	0.76
1:B:359:ARG:HG2	1:B:360:PRO:HD2	1.67	0.76
1:C:307:ASN:ND2	1:C:308:PRO:HD2	2.01	0.76
1:C:212:LEU:HD12	1:C:300:ILE:CD1	2.15	0.75
1:G:82:SER:HB2	1:G:189:SER:OG	1.86	0.75
1:C:189:SER:O	1:C:383:ILE:HG12	1.86	0.75
1:A:22:ILE:HG23	1:A:117:ILE:HD11	1.69	0.75
1:A:93:ARG:HA	1:A:108:HIS:CD2	2.22	0.75
1:B:51:ARG:HG3	1:B:51:ARG:HH11	1.51	0.75
1:F:212:LEU:HD12	1:F:300:ILE:HD12	1.69	0.75
1:H:38:GLU:HG2	1:H:88:ASP:HB3	1.69	0.74
1:G:212:LEU:HD12	1:G:300:ILE:CD1	2.18	0.74
1:E:49:MET:HB3	1:E:56:PRO:HG3	1.68	0.74
1:B:61:VAL:HG13	1:H:173:LEU:CD1	2.14	0.74
1:E:82:SER:HB2	1:E:189:SER:OG	1.88	0.74
1:A:116:ARG:HD3	1:A:400:TYR:CD2	2.23	0.73
1:D:173:LEU:CD1	1:F:61:VAL:HG13	2.17	0.73
1:A:114:GLU:OE1	1:A:116:ARG:HG2	1.88	0.73
1:H:212:LEU:HD12	1:H:300:ILE:HD12	1.71	0.73
1:A:6:ASP:HA	1:A:9:LYS:HG3	1.70	0.73
1:F:89:THR:HG21	1:F:121:SER:HB2	1.68	0.73
1:B:212:LEU:HD12	1:B:300:ILE:HD12	1.70	0.73
1:A:212:LEU:HD12	1:A:300:ILE:CD1	2.18	0.73
1:B:175:GLU:OE1	1:B:334:ARG:NE	2.23	0.72
1:F:51:ARG:HG3	1:F:51:ARG:HH11	1.55	0.72
1:D:212:LEU:HD12	1:D:300:ILE:HD12	1.72	0.72
1:B:212:LEU:HD22	1:B:328:VAL:CG1	2.19	0.72
1:D:212:LEU:HD22	1:D:328:VAL:CG1	2.20	0.72
1:A:114:GLU:O	1:A:116:ARG:N	2.23	0.72
1:F:343:GLU:HB3	1:F:344:PRO:HD3	1.72	0.72
1:D:61:VAL:HG13	1:F:173:LEU:HD11	1.72	0.71
1:F:221:SER:O	1:F:222:ARG:NH1	2.24	0.71
1:G:7:VAL:O	1:G:11:MET:HG2	1.91	0.71
1:B:398:ASN:OD1	1:B:399:THR:N	2.24	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:114:GLU:O	1:G:116:ARG:N	2.24	0.71
1:D:114:GLU:OE1	1:D:116:ARG:HG2	1.91	0.71
1:G:203:ALA:HB2	1:G:364:PRO:HG3	1.72	0.71
1:G:307:ASN:ND2	1:G:308:PRO:HD2	2.05	0.70
1:C:52:ASN:OD1	1:C:137:LYS:NZ	2.23	0.70
1:F:48:TRP:O	1:F:52:ASN:ND2	2.25	0.70
1:A:89:THR:HG21	1:A:121:SER:HB2	1.73	0.70
1:E:251:GLU:OE1	1:F:229:ARG:NH2	2.25	0.70
1:H:116:ARG:HD3	1:H:400:TYR:CD2	2.26	0.70
1:D:373:ARG:NH1	2:D:501:6R6:O12	2.24	0.69
1:E:307:ASN:ND2	1:E:308:PRO:HD2	2.05	0.69
1:A:173:LEU:CD1	1:C:61:VAL:HG13	2.22	0.69
1:F:279:ILE:HD12	1:F:296:ILE:HG22	1.72	0.69
1:H:48:TRP:CZ3	1:H:133:LEU:HD22	2.28	0.69
1:F:224:THR:HB	1:F:225:PRO:HD3	1.75	0.69
1:D:48:TRP:O	1:D:52:ASN:ND2	2.22	0.69
1:B:116:ARG:HD3	1:B:400:TYR:CD2	2.28	0.69
1:C:51:ARG:HG3	1:C:51:ARG:HH11	1.57	0.69
1:H:6:ASP:HA	1:H:9:LYS:HG3	1.74	0.69
1:G:43:ASP:OD1	1:G:58:ARG:NH1	2.26	0.69
1:A:38:GLU:CG	1:A:88:ASP:HB3	2.22	0.69
1:F:114:GLU:OE1	1:F:116:ARG:HG2	1.93	0.69
1:B:114:GLU:O	1:B:116:ARG:N	2.27	0.68
1:F:162:GLU:HG2	1:F:332:LEU:HD22	1.73	0.68
1:E:173:LEU:CD1	1:G:61:VAL:HG13	2.22	0.68
1:E:212:LEU:HD12	1:E:300:ILE:HD12	1.75	0.68
1:H:271:VAL:HG23	1:H:366:SER:HB3	1.73	0.68
1:A:373:ARG:NH1	2:A:501:6R6:O11	2.27	0.68
1:G:162:GLU:HG2	1:G:332:LEU:HD22	1.74	0.68
1:G:398:ASN:OD1	1:G:399:THR:N	2.27	0.68
1:G:373:ARG:NH1	2:G:501:6R6:O11	2.27	0.68
1:E:373:ARG:NH1	2:E:501:6R6:O11	2.27	0.68
1:D:43:ASP:OD1	1:D:58:ARG:NH1	2.27	0.68
1:C:116:ARG:HD3	1:C:400:TYR:CD2	2.29	0.67
1:E:114:GLU:OE1	1:E:116:ARG:HG2	1.94	0.67
1:G:116:ARG:HD3	1:G:400:TYR:CD2	2.28	0.67
1:C:212:LEU:HD22	1:C:328:VAL:CG1	2.25	0.67
1:E:22:ILE:HG23	1:E:117:ILE:HD11	1.76	0.67
1:C:266:GLY:O	1:C:363:ARG:NH1	2.27	0.67
1:D:243:MET:HE3	1:D:296:ILE:HG23	1.76	0.66
1:C:359:ARG:HG2	1:C:360:PRO:HD2	1.76	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:22:ILE:HG23	1:H:117:ILE:HD11	1.77	0.66
1:C:229:ARG:NH1	1:H:274:VAL:O	2.28	0.66
1:B:243:MET:HE3	1:B:296:ILE:HG23	1.77	0.66
1:E:374:ASP:C	1:E:377:PRO:HD2	2.16	0.66
1:H:342:ILE:HD11	1:H:386:LEU:HD23	1.77	0.66
1:B:91:MET:HG3	1:B:107:TYR:HB3	1.77	0.66
1:G:343:GLU:HB3	1:G:344:PRO:HD3	1.78	0.66
1:A:207:ALA:HB1	1:A:302:LEU:O	1.96	0.66
1:F:48:TRP:CZ3	1:F:133:LEU:HD22	2.30	0.66
1:D:7:VAL:O	1:D:11:MET:HG2	1.96	0.66
1:E:343:GLU:HB3	1:E:344:PRO:HD3	1.77	0.66
1:G:245:LYS:O	1:G:248:GLU:HG2	1.95	0.66
1:A:22:ILE:O	1:A:26:VAL:HG23	1.97	0.65
1:C:243:MET:HE3	1:C:296:ILE:HG23	1.78	0.65
1:D:51:ARG:HG3	1:D:51:ARG:HH11	1.59	0.65
1:G:93:ARG:HA	1:G:108:HIS:CD2	2.30	0.65
1:D:216:VAL:HG12	1:D:243:MET:CE	2.27	0.65
1:F:256:ASN:O	1:F:260:ARG:HG3	1.97	0.65
1:C:203:ALA:HB2	1:C:364:PRO:HG3	1.78	0.65
1:H:6:ASP:HA	1:H:9:LYS:HE2	1.79	0.65
1:H:48:TRP:O	1:H:52:ASN:ND2	2.24	0.65
1:H:51:ARG:HH11	1:H:51:ARG:HG3	1.61	0.65
1:A:374:ASP:C	1:A:377:PRO:HD2	2.17	0.65
1:C:48:TRP:O	1:C:52:ASN:ND2	2.29	0.65
1:G:52:ASN:OD1	1:G:137:LYS:NZ	2.30	0.65
1:A:51:ARG:HH11	1:A:51:ARG:HG3	1.61	0.65
1:C:105:ARG:O	1:C:109:GLU:HG3	1.97	0.65
1:D:7:VAL:HG22	1:D:418:MET:HE3	1.80	0.64
1:H:89:THR:HG21	1:H:121:SER:HB2	1.78	0.64
1:H:367:PRO:O	1:H:371:MET:HE3	1.98	0.64
1:G:212:LEU:HD12	1:G:300:ILE:HD12	1.78	0.64
1:H:7:VAL:HG22	1:H:418:MET:CE	2.26	0.64
1:C:243:MET:O	1:C:247:VAL:HG23	1.98	0.63
1:H:114:GLU:OE1	1:H:116:ARG:HG2	1.97	0.63
1:H:212:LEU:HD22	1:H:328:VAL:CG1	2.28	0.63
1:A:160:ASP:N	1:A:160:ASP:OD1	2.32	0.63
1:D:38:GLU:HG2	1:D:88:ASP:HB3	1.79	0.63
1:F:114:GLU:C	1:F:116:ARG:H	2.02	0.63
1:H:245:LYS:O	1:H:248:GLU:HG2	1.99	0.63
1:G:114:GLU:C	1:G:116:ARG:H	2.02	0.63
1:B:216:VAL:HG12	1:B:243:MET:CE	2.28	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:51:ARG:HH11	1:E:51:ARG:HG3	1.64	0.62
1:E:91:MET:HE3	1:E:120:TYR:CG	2.34	0.62
1:F:243:MET:O	1:F:247:VAL:HG23	2.00	0.62
1:A:48:TRP:CZ3	1:A:133:LEU:HD22	2.34	0.62
1:F:93:ARG:HA	1:F:108:HIS:CD2	2.35	0.62
1:H:52:ASN:OD1	1:H:137:LYS:NZ	2.33	0.62
1:B:213:LYS:HG3	1:B:297:TYR:CE2	2.34	0.62
1:B:43:ASP:OD1	1:B:58:ARG:NH1	2.33	0.62
1:A:229:ARG:NH1	1:G:274:VAL:O	2.33	0.62
1:D:114:GLU:C	1:D:116:ARG:H	2.02	0.62
1:E:38:GLU:HG2	1:E:88:ASP:HB3	1.82	0.62
1:F:396:GLY:N	1:F:397:GLY:HA2	2.14	0.62
1:F:75:THR:HG23	1:F:146:LEU:O	1.99	0.62
1:G:342:ILE:HD11	1:G:386:LEU:HD23	1.80	0.62
1:D:279:ILE:HD12	1:D:296:ILE:HG22	1.81	0.62
1:E:221:SER:O	1:E:222:ARG:NH1	2.33	0.62
1:B:245:LYS:O	1:B:248:GLU:HG2	2.00	0.62
1:G:350:GLU:O	1:G:354:ARG:HG3	1.99	0.61
1:A:118:TYR:CD1	1:A:400:TYR:HB3	2.35	0.61
1:G:89:THR:HG21	1:G:121:SER:HB2	1.82	0.61
1:A:43:ASP:OD1	1:A:58:ARG:NH1	2.33	0.61
1:C:6:ASP:HA	1:C:9:LYS:NZ	2.16	0.61
1:H:398:ASN:OD1	1:H:399:THR:N	2.33	0.61
1:A:7:VAL:HG22	1:A:418:MET:CE	2.30	0.61
1:A:75:THR:HG23	1:A:146:LEU:O	2.00	0.61
1:F:212:LEU:HD12	1:F:300:ILE:CD1	2.29	0.61
1:F:398:ASN:OD1	1:F:399:THR:N	2.34	0.61
1:A:86:HIS:CB	1:A:158:GLU:HB2	2.31	0.61
1:D:38:GLU:CG	1:D:88:ASP:HB3	2.31	0.61
1:G:212:LEU:HD22	1:G:328:VAL:CG1	2.30	0.61
1:G:6:ASP:OD1	1:G:9:LYS:HE2	2.00	0.61
1:A:243:MET:HE3	1:A:296:ILE:HG23	1.83	0.61
1:C:124:ASN:HA	1:C:125:CYS:HB2	1.83	0.61
1:A:274:VAL:O	1:G:229:ARG:NH1	2.33	0.60
1:H:22:ILE:O	1:H:26:VAL:HG23	2.01	0.60
1:E:114:GLU:C	1:E:116:ARG:H	2.03	0.60
1:F:374:ASP:C	1:F:377:PRO:HD2	2.22	0.60
1:A:279:ILE:HD12	1:A:296:ILE:HG22	1.83	0.60
1:D:48:TRP:CZ3	1:D:133:LEU:HD22	2.35	0.60
1:E:274:VAL:O	1:F:229:ARG:NH1	2.34	0.60
1:F:7:VAL:O	1:F:11:MET:HG2	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:160:ASP:N	1:B:160:ASP:OD1	2.34	0.60
1:F:22:ILE:HG23	1:F:117:ILE:CD1	2.31	0.60
1:G:160:ASP:OD1	1:G:160:ASP:N	2.33	0.60
1:H:38:GLU:CG	1:H:88:ASP:HB3	2.30	0.60
1:D:396:GLY:N	1:D:397:GLY:HA2	2.17	0.60
1:C:114:GLU:O	1:C:116:ARG:N	2.34	0.60
1:A:245:LYS:O	1:A:248:GLU:HG2	2.02	0.60
1:A:6:ASP:OD1	1:A:9:LYS:HE2	2.02	0.60
1:G:51:ARG:HG3	1:G:51:ARG:HH11	1.64	0.60
1:C:115:GLY:O	1:C:403:VAL:HG23	2.02	0.60
1:H:191:TYR:OH	1:H:344:PRO:HG2	2.02	0.60
1:D:205:VAL:HG13	1:D:369:CYS:HB3	1.84	0.59
1:D:61:VAL:HG13	1:F:173:LEU:HD13	1.83	0.59
1:E:93:ARG:HA	1:E:108:HIS:CD2	2.36	0.59
1:A:256:ASN:O	1:A:260:ARG:HG3	2.02	0.59
1:C:245:LYS:O	1:C:248:GLU:HG2	2.01	0.59
1:D:49:MET:HB3	1:D:56:PRO:HG3	1.84	0.59
1:E:396:GLY:N	1:E:397:GLY:HA2	2.16	0.59
1:G:22:ILE:O	1:G:26:VAL:HG23	2.02	0.59
1:D:106:ILE:HD11	1:D:399:THR:HB	1.83	0.59
1:G:364:PRO:HG2	1:G:369:CYS:SG	2.42	0.59
1:H:6:ASP:OD1	1:H:9:LYS:HE2	2.03	0.59
1:A:7:VAL:HA	1:A:418:MET:HE1	1.84	0.59
1:F:212:LEU:HD22	1:F:328:VAL:CG1	2.33	0.59
1:A:24:THR:O	1:A:28:LEU:HB2	2.02	0.59
1:B:114:GLU:C	1:B:116:ARG:H	2.06	0.59
1:B:89:THR:HG21	1:B:121:SER:HB2	1.84	0.59
1:D:6:ASP:HA	1:D:9:LYS:HE2	1.84	0.59
1:F:414:ALA:O	1:F:418:MET:HG3	2.03	0.59
1:A:86:HIS:HB3	1:A:158:GLU:HB2	1.85	0.59
1:A:265:TYR:OH	1:A:305:ASP:OD2	2.16	0.59
1:B:18:ARG:HD2	1:B:403:VAL:HG12	1.84	0.59
1:A:57:GLU:OE1	1:C:329:LYS:NZ	2.36	0.58
1:F:52:ASN:OD1	1:F:137:LYS:NZ	2.35	0.58
1:H:17:MET:CE	1:H:137:LYS:HD3	2.33	0.58
1:H:374:ASP:O	1:H:377:PRO:HD2	2.03	0.58
1:A:114:GLU:C	1:A:116:ARG:H	2.06	0.58
1:H:114:GLU:C	1:H:116:ARG:H	2.06	0.58
1:E:75:THR:HG23	1:E:146:LEU:O	2.03	0.58
1:E:204:TRP:CE2	1:E:267:GLY:HA2	2.39	0.58
1:G:48:TRP:CZ3	1:G:133:LEU:HD22	2.38	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:160:ASP:OD1	1:F:160:ASP:N	2.36	0.58
1:G:224:THR:HB	1:G:225:PRO:HD3	1.85	0.58
1:A:83:PHE:HE1	1:A:193:LEU:HD22	1.69	0.58
1:B:243:MET:O	1:B:247:VAL:HG23	2.03	0.58
1:B:93:ARG:HA	1:B:108:HIS:CD2	2.38	0.58
1:H:396:GLY:N	1:H:397:GLY:HA2	2.18	0.58
1:H:374:ASP:C	1:H:377:PRO:HD2	2.24	0.58
1:F:116:ARG:HD3	1:F:400:TYR:CD2	2.39	0.58
1:G:189:SER:O	1:G:383:ILE:HG12	2.04	0.58
1:H:359:ARG:HG2	1:H:360:PRO:HD2	1.85	0.58
1:C:11:MET:SD	1:C:411:LYS:HG2	2.44	0.57
1:E:160:ASP:OD1	1:E:160:ASP:N	2.37	0.57
1:G:261:TYR:OH	1:G:314:GLU:OE2	2.15	0.57
1:G:396:GLY:N	1:G:397:GLY:HA2	2.19	0.57
1:H:7:VAL:O	1:H:11:MET:HG2	2.04	0.57
1:F:364:PRO:HG2	1:F:369:CYS:SG	2.43	0.57
1:H:91:MET:HE3	1:H:120:TYR:CG	2.39	0.57
1:H:418:MET:O	1:H:422:ASN:ND2	2.37	0.57
1:B:51:ARG:HH11	1:B:51:ARG:CG	2.18	0.57
1:E:329:LYS:NZ	1:G:57:GLU:OE1	2.38	0.57
1:B:216:VAL:HG12	1:B:243:MET:HE1	1.86	0.57
1:G:77:GLY:O	1:G:424:THR:HG22	2.04	0.57
1:E:224:THR:HB	1:E:225:PRO:HD3	1.85	0.57
1:E:224:THR:O	1:E:227:VAL:HG12	2.05	0.57
1:E:374:ASP:O	1:E:377:PRO:HD2	2.03	0.57
1:F:18:ARG:HD3	1:F:407:LEU:HD22	1.87	0.57
1:H:160:ASP:N	1:H:160:ASP:OD1	2.37	0.57
1:B:88:ASP:OD2	1:B:159:ILE:HG13	2.04	0.57
1:C:160:ASP:OD1	1:C:160:ASP:N	2.37	0.57
1:C:22:ILE:HG23	1:C:117:ILE:CD1	2.29	0.57
1:G:75:THR:HG23	1:G:146:LEU:O	2.04	0.57
1:G:359:ARG:HG2	1:G:360:PRO:HD2	1.86	0.57
1:G:374:ASP:C	1:G:377:PRO:HD2	2.26	0.57
1:B:314:GLU:O	1:B:317:ALA:HB3	2.05	0.57
1:E:212:LEU:HD22	1:E:328:VAL:CG1	2.35	0.57
1:A:307:ASN:ND2	1:A:308:PRO:HD2	2.19	0.57
1:A:343:GLU:HB3	1:A:344:PRO:HD3	1.87	0.57
1:D:343:GLU:HB3	1:D:344:PRO:HD3	1.86	0.57
1:H:285:TYR:CD1	1:H:286:LYS:HG3	2.39	0.57
1:A:419:ASP:OD1	1:A:423:ARG:NE	2.34	0.56
1:F:353:HIS:CE1	1:F:361:THR:HG22	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:93:ARG:HA	1:H:108:HIS:CD2	2.40	0.56
1:D:374:ASP:O	1:D:377:PRO:HD2	2.05	0.56
1:F:201:LYS:HD2	1:F:364:PRO:HA	1.88	0.56
1:F:203:ALA:CB	1:F:364:PRO:HG3	2.30	0.56
1:F:38:GLU:HG2	1:F:88:ASP:HB3	1.86	0.56
1:B:396:GLY:N	1:B:397:GLY:HA2	2.20	0.56
1:C:279:ILE:HD12	1:C:296:ILE:HG22	1.87	0.56
1:C:374:ASP:C	1:C:377:PRO:HD2	2.26	0.56
1:F:51:ARG:CG	1:F:51:ARG:HH11	2.18	0.56
1:B:191:TYR:OH	1:B:344:PRO:HG2	2.06	0.56
1:B:75:THR:HG23	1:B:146:LEU:O	2.05	0.56
1:D:285:TYR:CD1	1:D:286:LYS:HG3	2.41	0.56
1:E:43:ASP:OD1	1:E:58:ARG:NH1	2.38	0.56
1:G:86:HIS:CB	1:G:158:GLU:HB2	2.36	0.56
1:A:221:SER:O	1:A:222:ARG:NH1	2.38	0.56
1:B:38:GLU:HG2	1:B:88:ASP:HB3	1.87	0.56
1:E:132:TRP:HZ3	1:E:152:LEU:HB3	1.71	0.56
1:G:18:ARG:HD2	1:G:403:VAL:HG12	1.87	0.56
1:H:216:VAL:HG12	1:H:243:MET:HE1	1.88	0.56
1:A:42:GLY:HA2	1:A:87:LEU:HD13	1.87	0.56
1:E:309:LEU:HD21	1:G:188:ILE:HG13	1.88	0.56
1:B:274:VAL:O	1:D:229:ARG:NH1	2.39	0.56
1:B:287:ILE:HB	1:D:277:GLY:HA3	1.88	0.56
1:H:105:ARG:O	1:H:109:GLU:HG3	2.06	0.56
1:H:216:VAL:HG12	1:H:243:MET:CE	2.36	0.56
1:H:243:MET:HE3	1:H:296:ILE:HG23	1.87	0.56
1:C:342:ILE:HD11	1:C:386:LEU:HD23	1.87	0.56
1:H:106:ILE:HG23	1:H:107:TYR:CD1	2.40	0.56
1:A:398:ASN:OD1	1:A:399:THR:N	2.39	0.55
1:D:213:LYS:HG3	1:D:297:TYR:CE2	2.41	0.55
1:F:374:ASP:O	1:F:377:PRO:HD2	2.06	0.55
1:A:213:LYS:HG3	1:A:297:TYR:CE2	2.42	0.55
1:A:7:VAL:O	1:A:11:MET:HG2	2.06	0.55
1:A:418:MET:O	1:A:422:ASN:ND2	2.38	0.55
1:B:285:TYR:CD1	1:B:286:LYS:HG3	2.42	0.55
1:A:212:LEU:HD22	1:A:328:VAL:CG1	2.37	0.55
1:B:332:LEU:HD12	1:H:184:SER:O	2.05	0.55
1:C:18:ARG:O	1:C:22:ILE:HG13	2.07	0.55
1:D:7:VAL:HA	1:D:418:MET:HE1	1.86	0.55
1:B:18:ARG:HD2	1:B:403:VAL:CG1	2.37	0.55
1:B:373:ARG:NH1	2:B:501:6R6:O11	2.40	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:11:MET:SD	1:D:411:LYS:HG2	2.47	0.55
1:D:93:ARG:HA	1:D:108:HIS:CD2	2.41	0.55
1:D:82:SER:HB2	1:D:189:SER:HG	1.72	0.55
1:H:76:GLY:O	1:H:147:LYS:HB3	2.07	0.55
1:B:364:PRO:HG2	1:B:369:CYS:SG	2.47	0.54
1:C:38:GLU:HG2	1:C:88:ASP:HB3	1.88	0.54
1:C:114:GLU:C	1:C:116:ARG:H	2.10	0.54
1:E:7:VAL:O	1:E:11:MET:HG2	2.08	0.54
1:F:63:ASP:N	1:F:63:ASP:OD1	2.23	0.54
1:G:204:TRP:CE2	1:G:267:GLY:HA2	2.42	0.54
1:H:63:ASP:OD1	1:H:63:ASP:N	2.39	0.54
1:E:22:ILE:HG23	1:E:117:ILE:CD1	2.38	0.54
1:C:396:GLY:N	1:C:397:GLY:HA2	2.23	0.54
1:C:93:ARG:HA	1:C:108:HIS:CD2	2.42	0.54
1:E:319:VAL:HG11	1:E:326:ALA:HB3	1.88	0.54
1:G:314:GLU:O	1:G:317:ALA:HB3	2.08	0.54
1:G:262:THR:HG23	1:G:271:VAL:HG22	1.88	0.54
1:G:83:PHE:HE1	1:G:193:LEU:HD22	1.72	0.54
1:C:224:THR:HB	1:C:225:PRO:HD3	1.89	0.54
1:C:354:ARG:NH1	1:C:360:PRO:HG3	2.23	0.54
1:H:14:LEU:HD23	1:H:138:ALA:HB2	1.89	0.54
1:H:256:ASN:O	1:H:260:ARG:HG3	2.08	0.54
1:B:105:ARG:O	1:B:109:GLU:HG3	2.08	0.54
1:C:374:ASP:O	1:C:377:PRO:HD2	2.07	0.54
1:D:207:ALA:HB1	1:D:302:LEU:O	2.08	0.54
1:B:224:THR:HB	1:B:225:PRO:HD3	1.89	0.54
1:B:115:GLY:O	1:B:403:VAL:HG23	2.08	0.54
1:B:204:TRP:CE2	1:B:267:GLY:HA2	2.43	0.54
1:B:81:LEU:HD12	1:B:82:SER:H	1.73	0.54
1:F:314:GLU:O	1:F:317:ALA:HB3	2.08	0.54
1:D:256:ASN:O	1:D:260:ARG:HG3	2.08	0.54
1:H:240:ILE:HD12	1:H:291:PRO:HG2	1.90	0.54
1:H:240:ILE:CD1	1:H:291:PRO:HG2	2.38	0.54
1:F:82:SER:HB2	1:F:189:SER:HG	1.73	0.53
1:A:6:ASP:HA	1:A:9:LYS:CE	2.35	0.53
1:D:398:ASN:OD1	1:D:399:THR:N	2.41	0.53
1:F:204:TRP:CE2	1:F:267:GLY:HA2	2.43	0.53
1:F:279:ILE:HD12	1:F:296:ILE:CG2	2.39	0.53
1:F:7:VAL:HG22	1:F:418:MET:HE3	1.89	0.53
1:D:216:VAL:HG12	1:D:243:MET:HE1	1.89	0.53
1:D:262:THR:HG23	1:D:271:VAL:HG22	1.88	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:7:VAL:HG22	1:D:418:MET:CE	2.38	0.53
1:D:86:HIS:CB	1:D:158:GLU:HB2	2.38	0.53
1:F:107:TYR:HB2	1:F:108:HIS:CE1	2.44	0.53
1:B:343:GLU:HB3	1:B:344:PRO:HD3	1.90	0.53
1:H:86:HIS:CD2	1:H:86:HIS:H	2.26	0.53
1:E:89:THR:HG21	1:E:121:SER:HB2	1.91	0.53
1:H:51:ARG:HH11	1:H:51:ARG:CG	2.22	0.53
1:A:396:GLY:N	1:A:397:GLY:HA2	2.21	0.53
1:F:245:LYS:O	1:F:248:GLU:HG2	2.08	0.53
1:C:274:VAL:O	1:H:229:ARG:NH1	2.42	0.53
1:A:6:ASP:O	1:A:9:LYS:HG3	2.08	0.53
1:D:89:THR:HG21	1:D:121:SER:HB2	1.90	0.53
1:E:129:MET:HA	1:E:132:TRP:NE1	2.24	0.53
1:E:398:ASN:OD1	1:E:399:THR:N	2.42	0.53
1:F:6:ASP:HA	1:F:9:LYS:HE2	1.91	0.53
1:G:51:ARG:CG	1:G:51:ARG:HH11	2.22	0.53
1:A:359:ARG:HG2	1:A:360:PRO:HD2	1.90	0.53
1:B:22:ILE:O	1:B:26:VAL:HG23	2.09	0.53
1:D:212:LEU:HD22	1:D:328:VAL:HG12	1.89	0.53
1:D:51:ARG:CG	1:D:51:ARG:HH11	2.22	0.53
1:F:262:THR:HG23	1:F:271:VAL:HG22	1.91	0.53
1:G:256:ASN:O	1:G:260:ARG:HG3	2.09	0.53
1:C:373:ARG:NH1	2:C:501:6R6:O12	2.42	0.52
1:D:342:ILE:HD11	1:D:386:LEU:HD23	1.92	0.52
1:E:106:ILE:HD11	1:E:399:THR:HB	1.90	0.52
1:F:105:ARG:O	1:F:109:GLU:HG3	2.09	0.52
2:A:501:6R6:O09	1:G:223:TYR:HB2	2.09	0.52
1:B:221:SER:O	1:B:222:ARG:NH1	2.42	0.52
1:C:285:TYR:CD1	1:C:286:LYS:HG3	2.45	0.52
1:E:38:GLU:CG	1:E:88:ASP:HB3	2.39	0.52
1:G:238:ASN:OD1	1:G:240:ILE:N	2.39	0.52
1:A:93:ARG:HA	1:A:108:HIS:HD2	1.72	0.52
1:B:256:ASN:O	1:B:260:ARG:HG3	2.10	0.52
1:D:374:ASP:C	1:D:377:PRO:HD2	2.30	0.52
1:H:110:ALA:HA	1:H:118:TYR:O	2.09	0.52
1:F:213:LYS:HG3	1:F:297:TYR:CE2	2.45	0.52
1:G:353:HIS:CE1	1:G:361:THR:HG22	2.44	0.52
1:B:162:GLU:HG2	1:B:332:LEU:HD22	1.91	0.52
1:C:211:PHE:HB2	1:C:332:LEU:HB3	1.91	0.52
1:C:270:VAL:HG11	1:C:303:ASN:HB3	1.91	0.52
1:D:162:GLU:HG2	1:D:332:LEU:HD22	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:352:ALA:O	1:A:356:VAL:HG23	2.09	0.52
1:D:6:ASP:O	1:D:9:LYS:HG3	2.08	0.52
1:B:201:LYS:HD2	1:B:364:PRO:HA	1.92	0.52
1:E:229:ARG:NH1	1:F:274:VAL:O	2.43	0.52
1:C:262:THR:HG23	1:C:271:VAL:HG22	1.91	0.52
1:G:88:ASP:OD1	1:G:88:ASP:N	2.43	0.52
1:H:366:SER:O	1:H:370:SER:OG	2.21	0.52
1:E:260:ARG:HH22	1:E:314:GLU:CD	2.13	0.51
1:E:391:GLY:HA3	1:E:401:PHE:CE1	2.45	0.51
1:A:23:GLN:O	1:A:27:GLU:HG3	2.10	0.51
1:A:116:ARG:HD3	1:A:400:TYR:CE2	2.44	0.51
1:C:22:ILE:O	1:C:26:VAL:HG23	2.10	0.51
1:C:51:ARG:CG	1:C:51:ARG:HH11	2.21	0.51
1:E:243:MET:O	1:E:247:VAL:HG23	2.10	0.51
1:C:221:SER:O	1:C:222:ARG:NH1	2.44	0.51
1:D:106:ILE:HG23	1:D:107:TYR:CD1	2.46	0.51
1:D:204:TRP:CE2	1:D:267:GLY:HA2	2.46	0.51
1:E:277:GLY:HA3	1:F:287:ILE:HB	1.92	0.51
1:B:374:ASP:C	1:B:377:PRO:HD2	2.31	0.51
1:C:77:GLY:O	1:C:424:THR:HG22	2.11	0.51
1:E:203:ALA:HB2	1:E:364:PRO:HG3	1.93	0.51
1:H:43:ASP:OD1	1:H:58:ARG:NH1	2.42	0.51
1:H:213:LYS:HG3	1:H:297:TYR:CE2	2.46	0.51
1:G:221:SER:O	1:G:222:ARG:NH1	2.44	0.51
1:G:7:VAL:HG22	1:G:418:MET:HE3	1.93	0.51
1:B:26:VAL:O	1:B:30:SER:OG	2.26	0.51
1:C:137:LYS:HE3	1:C:141:GLU:OE2	2.10	0.51
1:H:343:GLU:HB3	1:H:344:PRO:HD3	1.92	0.51
1:H:86:HIS:CB	1:H:158:GLU:HB2	2.40	0.51
1:B:342:ILE:HD11	1:B:345:LEU:HD23	1.93	0.50
1:D:300:ILE:HG21	1:D:311:VAL:HG11	1.93	0.50
1:A:229:ARG:NH2	1:G:251:GLU:OE1	2.44	0.50
1:B:132:TRP:HZ3	1:B:152:LEU:HB3	1.76	0.50
1:E:245:LYS:O	1:E:248:GLU:HG2	2.10	0.50
1:G:191:TYR:OH	1:G:344:PRO:HG2	2.10	0.50
1:E:23:GLN:O	1:E:27:GLU:HG3	2.11	0.50
1:D:184:SER:O	1:F:332:LEU:HD12	2.11	0.50
1:B:157:GLY:HA3	1:B:174:ALA:O	2.11	0.50
1:C:59:VAL:O	1:C:67:ASN:ND2	2.44	0.50
1:E:86:HIS:CB	1:E:158:GLU:HB2	2.41	0.50
1:G:197:ALA:HB2	1:G:401:PHE:HZ	1.77	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:263:ARG:HB3	1:H:265:TYR:HE1	1.76	0.50
1:A:342:ILE:HD11	1:A:386:LEU:HD23	1.92	0.50
1:A:83:PHE:CE1	1:A:193:LEU:HD22	2.47	0.50
1:C:175:GLU:OE1	1:C:334:ARG:NE	2.45	0.50
1:C:197:ALA:HB2	1:C:401:PHE:HZ	1.76	0.50
1:E:105:ARG:O	1:E:109:GLU:HG3	2.11	0.50
1:E:211:PHE:O	1:E:330:PRO:HA	2.10	0.50
1:F:22:ILE:O	1:F:26:VAL:HG23	2.12	0.50
1:G:163:PRO:HG3	1:G:173:LEU:HD23	1.93	0.50
1:B:251:GLU:OE1	1:D:229:ARG:NH2	2.44	0.50
1:B:374:ASP:O	1:B:377:PRO:HD2	2.11	0.50
1:D:243:MET:O	1:D:247:VAL:HG23	2.12	0.50
1:E:184:SER:O	1:G:332:LEU:HD12	2.11	0.50
1:A:224:THR:HB	1:A:225:PRO:HD3	1.92	0.50
1:E:24:THR:O	1:E:28:LEU:HB2	2.10	0.50
1:F:6:ASP:OD1	1:F:9:LYS:HE2	2.12	0.50
1:G:91:MET:HE3	1:G:120:TYR:CG	2.46	0.50
1:G:285:TYR:CD1	1:G:286:LYS:HG3	2.46	0.50
1:A:277:GLY:HA3	1:G:287:ILE:HB	1.94	0.50
1:C:28:LEU:HD22	1:C:126:LYS:HD2	1.94	0.50
1:E:196:GLU:O	1:E:389:GLY:HA3	2.12	0.50
1:G:88:ASP:OD2	1:G:159:ILE:HG13	2.11	0.50
1:H:7:VAL:HA	1:H:418:MET:HE1	1.92	0.50
1:C:203:ALA:O	1:C:387:THR:HG22	2.12	0.50
1:E:118:TYR:CD1	1:E:400:TYR:HB3	2.46	0.50
1:H:75:THR:HG23	1:H:146:LEU:O	2.11	0.50
1:B:86:HIS:H	1:B:86:HIS:CD2	2.30	0.49
1:F:211:PHE:O	1:F:330:PRO:HA	2.11	0.49
1:G:118:TYR:CD1	1:G:400:TYR:HB3	2.47	0.49
1:G:243:MET:O	1:G:247:VAL:HG23	2.12	0.49
1:G:24:THR:O	1:G:28:LEU:HB2	2.12	0.49
1:A:110:ALA:HA	1:A:118:TYR:O	2.12	0.49
1:A:51:ARG:HH11	1:A:51:ARG:CG	2.25	0.49
1:B:77:GLY:O	1:B:424:THR:HG22	2.12	0.49
1:C:334:ARG:HB3	1:C:372:TRP:CZ3	2.47	0.49
1:H:162:GLU:HG2	1:H:332:LEU:HD13	1.93	0.49
1:A:22:ILE:HG23	1:A:117:ILE:CD1	2.41	0.49
1:C:207:ALA:HB1	1:C:302:LEU:O	2.12	0.49
1:E:189:SER:O	1:E:383:ILE:HG12	2.13	0.49
1:D:6:ASP:OD1	1:D:9:LYS:HE2	2.12	0.49
1:E:124:ASN:HA	1:E:125:CYS:HB2	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:213:LYS:HG3	1:E:297:TYR:CE2	2.48	0.49
1:E:51:ARG:HH11	1:E:51:ARG:CG	2.25	0.49
1:H:224:THR:O	1:H:227:VAL:HG12	2.13	0.49
1:B:197:ALA:HB2	1:B:401:PHE:HZ	1.77	0.49
1:D:160:ASP:N	1:D:160:ASP:OD1	2.45	0.49
1:B:211:PHE:HB2	1:B:332:LEU:HB3	1.95	0.49
1:C:88:ASP:OD2	1:C:159:ILE:HG13	2.13	0.49
1:A:184:SER:O	1:C:332:LEU:HD12	2.13	0.49
1:C:408:LYS:O	1:C:411:LYS:N	2.46	0.49
1:D:88:ASP:N	1:D:88:ASP:OD1	2.46	0.49
1:F:203:ALA:HB2	1:F:364:PRO:CG	2.34	0.49
1:F:285:TYR:CD1	1:F:286:LYS:HG3	2.48	0.49
1:G:386:LEU:HD12	1:G:386:LEU:C	2.33	0.49
1:G:81:LEU:HD12	1:G:82:SER:H	1.78	0.49
1:A:243:MET:O	1:A:247:VAL:HG23	2.13	0.49
1:A:63:ASP:N	1:A:63:ASP:OD1	2.33	0.49
1:E:106:ILE:HG23	1:E:107:TYR:CD1	2.48	0.49
1:E:6:ASP:OD1	1:E:9:LYS:HE2	2.13	0.49
1:F:86:HIS:CB	1:F:158:GLU:HB2	2.43	0.49
1:A:204:TRP:CE2	1:A:267:GLY:HA2	2.48	0.49
1:B:194:VAL:CG2	1:B:375:THR:HG22	2.43	0.49
1:C:6:ASP:O	1:C:9:LYS:HG3	2.12	0.49
1:D:212:LEU:CD1	1:D:300:ILE:CD1	2.87	0.49
1:D:271:VAL:HG23	1:D:366:SER:HB3	1.95	0.49
1:C:213:LYS:HG3	1:C:297:TYR:CE2	2.48	0.48
1:E:266:GLY:O	1:E:363:ARG:NH1	2.46	0.48
1:F:24:THR:O	1:F:28:LEU:HB2	2.12	0.48
1:G:86:HIS:HB3	1:G:158:GLU:HB2	1.94	0.48
1:E:229:ARG:NH2	1:F:251:GLU:OE1	2.46	0.48
1:F:224:THR:O	1:F:227:VAL:HG12	2.13	0.48
1:G:201:LYS:HD2	1:G:364:PRO:HA	1.95	0.48
1:H:24:THR:O	1:H:28:LEU:HB2	2.12	0.48
1:A:6:ASP:CA	1:A:9:LYS:HG3	2.41	0.48
1:B:18:ARG:O	1:B:22:ILE:HG13	2.13	0.48
1:C:212:LEU:CD1	1:C:300:ILE:CD1	2.89	0.48
1:D:132:TRP:HZ3	1:D:152:LEU:HB3	1.79	0.48
1:A:347:ASN:O	1:A:351:VAL:HG23	2.13	0.48
1:E:211:PHE:HB2	1:E:332:LEU:HB3	1.95	0.48
1:E:312:GLN:NE2	1:E:316:GLU:OE2	2.47	0.48
1:F:265:TYR:OH	1:F:305:ASP:OD2	2.17	0.48
1:E:256:ASN:O	1:E:260:ARG:HG3	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:129:MET:O	1:G:133:LEU:HB2	2.14	0.48
1:H:86:HIS:ND1	1:H:158:GLU:HB3	2.28	0.48
1:B:6:ASP:O	1:B:9:LYS:HG3	2.13	0.48
1:B:124:ASN:HA	1:B:125:CYS:HB2	1.94	0.48
1:D:216:VAL:HG12	1:D:243:MET:HE2	1.96	0.48
1:F:38:GLU:CG	1:F:88:ASP:HB3	2.44	0.48
1:A:76:GLY:O	1:A:147:LYS:HB3	2.13	0.48
1:A:203:ALA:HB2	1:A:364:PRO:HG3	1.96	0.48
1:E:216:VAL:HG12	1:E:243:MET:CE	2.44	0.48
1:G:38:GLU:HG2	1:G:88:ASP:HB3	1.96	0.48
1:A:11:MET:SD	1:A:411:LYS:HG2	2.53	0.47
1:D:245:LYS:HD2	1:D:245:LYS:HA	1.67	0.47
1:D:24:THR:O	1:D:28:LEU:HB2	2.14	0.47
1:E:206:GLU:HA	1:E:206:GLU:OE1	2.13	0.47
1:H:224:THR:HB	1:H:225:PRO:HD3	1.95	0.47
1:B:207:ALA:HB1	1:B:302:LEU:O	2.15	0.47
1:B:191:TYR:HD1	1:B:384:PRO:HB2	1.78	0.47
1:G:86:HIS:CD2	1:G:86:HIS:H	2.32	0.47
1:A:332:LEU:HD12	1:C:184:SER:O	2.15	0.47
1:D:343:GLU:OE2	1:D:347:ASN:HB2	2.15	0.47
1:H:195:ALA:O	1:H:196:GLU:HG2	2.14	0.47
1:A:271:VAL:HG23	1:A:366:SER:HB3	1.96	0.47
1:A:260:ARG:HH22	1:A:314:GLU:CD	2.18	0.47
1:B:11:MET:SD	1:B:411:LYS:HG2	2.54	0.47
1:G:23:GLN:O	1:G:27:GLU:HG3	2.14	0.47
1:D:180:ARG:NH2	1:F:380:GLU:OE1	2.48	0.47
1:D:83:PHE:CE1	1:D:193:LEU:HD22	2.49	0.47
1:E:11:MET:SD	1:E:411:LYS:HG2	2.54	0.47
1:H:200:PHE:HD1	1:H:388:TYR:HH	1.63	0.47
1:A:91:MET:CE	1:A:120:TYR:CG	2.97	0.47
1:D:189:SER:O	1:D:383:ILE:HG12	2.14	0.47
1:F:114:GLU:C	1:F:116:ARG:N	2.65	0.47
1:E:224:THR:OG1	1:F:299:ASP:OD2	2.26	0.47
1:G:213:LYS:HG3	1:G:297:TYR:CE2	2.49	0.47
1:A:216:VAL:HG12	1:A:243:MET:CE	2.44	0.47
1:E:81:LEU:HD12	1:E:82:SER:H	1.79	0.47
1:G:106:ILE:HG23	1:G:107:TYR:CD1	2.50	0.47
1:H:106:ILE:HD11	1:H:399:THR:HB	1.97	0.47
1:H:207:ALA:HB1	1:H:302:LEU:O	2.15	0.47
1:A:212:LEU:CD1	1:A:300:ILE:CD1	2.89	0.47
1:A:196:GLU:HB3	1:A:373:ARG:NH1	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:22:ILE:HG23	1:D:117:ILE:CD1	2.34	0.47
1:D:86:HIS:CD2	1:D:86:HIS:H	2.33	0.47
1:G:114:GLU:C	1:G:116:ARG:N	2.65	0.47
1:H:223:TYR:CD1	1:H:225:PRO:HD2	2.49	0.47
1:D:223:TYR:CG	1:D:225:PRO:HD2	2.50	0.47
1:H:48:TRP:CE3	1:H:133:LEU:HD22	2.49	0.47
1:C:204:TRP:CE2	1:C:267:GLY:HA2	2.49	0.47
1:D:240:ILE:HD11	1:D:291:PRO:CG	2.45	0.47
1:E:63:ASP:N	1:E:63:ASP:OD1	2.37	0.47
1:G:157:GLY:HA3	1:G:174:ALA:O	2.14	0.47
1:H:240:ILE:CD1	1:H:291:PRO:CG	2.93	0.47
1:E:240:ILE:HD11	1:E:291:PRO:CG	2.45	0.46
1:G:283:VAL:HA	1:G:284:PRO:HD3	1.77	0.46
1:H:364:PRO:HG2	1:H:369:CYS:SG	2.54	0.46
1:B:106:ILE:HG23	1:B:107:TYR:CD1	2.50	0.46
1:B:195:ALA:O	1:B:196:GLU:HG2	2.16	0.46
1:B:216:VAL:CG1	1:B:243:MET:CE	2.93	0.46
1:C:125:CYS:SG	1:C:125:CYS:O	2.73	0.46
1:C:240:ILE:HD12	1:C:291:PRO:CD	2.45	0.46
1:G:376:ASN:HB2	1:G:377:PRO:HD3	1.98	0.46
1:H:309:LEU:HD23	1:H:309:LEU:HA	1.56	0.46
1:A:240:ILE:CD1	1:A:291:PRO:HG2	2.45	0.46
1:C:47:GLU:O	1:C:51:ARG:HB2	2.14	0.46
1:E:139:LEU:HD21	1:E:414:ALA:HB1	1.97	0.46
1:E:86:HIS:CD2	1:E:86:HIS:H	2.33	0.46
1:G:22:ILE:HG23	1:G:117:ILE:HD13	1.96	0.46
1:H:245:LYS:HA	1:H:245:LYS:HD2	1.49	0.46
1:A:14:LEU:HD23	1:A:138:ALA:HB2	1.97	0.46
1:F:18:ARG:O	1:F:21:LEU:HB3	2.16	0.46
1:F:283:VAL:HA	1:F:284:PRO:HD3	1.69	0.46
1:F:81:LEU:HD12	1:F:82:SER:H	1.81	0.46
1:G:265:TYR:OH	1:G:305:ASP:OD2	2.16	0.46
1:A:285:TYR:CD1	1:A:286:LYS:HG3	2.51	0.46
1:C:334:ARG:HB3	1:C:372:TRP:HZ3	1.80	0.46
1:D:86:HIS:HB3	1:D:158:GLU:HB2	1.97	0.46
1:E:86:HIS:ND1	1:E:158:GLU:HB3	2.31	0.46
1:F:359:ARG:HG2	1:F:360:PRO:HD2	1.98	0.46
1:A:234:LEU:HD22	1:A:324:LEU:HD21	1.97	0.46
1:E:260:ARG:NH2	1:E:314:GLU:OE2	2.48	0.46
1:C:216:VAL:CG1	1:C:243:MET:HE2	2.46	0.46
1:C:314:GLU:O	1:C:317:ALA:HB3	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:245:LYS:HD2	1:F:245:LYS:HA	1.57	0.46
1:A:116:ARG:HD2	1:A:118:TYR:OH	2.15	0.46
1:B:63:ASP:N	1:B:63:ASP:OD1	2.35	0.46
1:E:359:ARG:HG2	1:E:360:PRO:HD2	1.98	0.46
1:F:386:LEU:HD12	1:F:386:LEU:C	2.36	0.46
1:H:14:LEU:CD1	1:H:411:LYS:HG3	2.46	0.46
1:A:114:GLU:C	1:A:116:ARG:N	2.69	0.46
1:C:105:ARG:HG2	1:C:111:TRP:CZ3	2.50	0.46
1:C:270:VAL:HG11	1:C:303:ASN:CB	2.46	0.46
1:F:11:MET:SD	1:F:411:LYS:HG2	2.57	0.46
1:A:251:GLU:OE1	1:G:229:ARG:NH2	2.49	0.46
1:G:270:VAL:HG22	1:G:369:CYS:O	2.16	0.46
1:A:86:HIS:CD2	1:A:86:HIS:H	2.34	0.45
1:B:216:VAL:CG1	1:B:243:MET:HE2	2.46	0.45
1:C:391:GLY:HA3	1:C:401:PHE:CE1	2.51	0.45
1:F:125:CYS:SG	1:F:125:CYS:O	2.73	0.45
1:G:83:PHE:CE1	1:G:193:LEU:HD22	2.50	0.45
1:H:17:MET:HE2	1:H:137:LYS:HD3	1.97	0.45
1:H:279:ILE:HD12	1:H:296:ILE:HG22	1.98	0.45
1:A:250:LEU:HD23	1:A:318:VAL:HG11	1.98	0.45
1:D:206:GLU:HA	1:D:206:GLU:OE1	2.17	0.45
1:D:332:LEU:HD12	1:F:184:SER:O	2.16	0.45
1:E:245:LYS:HD2	1:E:245:LYS:HA	1.70	0.45
1:E:332:LEU:HD12	1:G:184:SER:O	2.16	0.45
1:G:93:ARG:HA	1:G:108:HIS:HD2	1.80	0.45
1:A:12:LYS:HA	1:A:12:LYS:HD3	1.70	0.45
1:A:34:PRO:HB3	1:A:92:ALA:HA	1.98	0.45
1:F:350:GLU:O	1:F:354:ARG:HG3	2.15	0.45
1:F:86:HIS:CD2	1:F:86:HIS:H	2.33	0.45
1:H:366:SER:N	1:H:367:PRO:CD	2.79	0.45
1:B:380:GLU:OE1	1:H:180:ARG:NH2	2.50	0.45
1:C:24:THR:O	1:C:28:LEU:HB2	2.16	0.45
1:D:195:ALA:O	1:D:196:GLU:HG2	2.16	0.45
1:D:63:ASP:OD1	1:D:63:ASP:N	2.30	0.45
1:E:238:ASN:ND2	1:E:290:PHE:HB2	2.31	0.45
1:E:332:LEU:HD12	1:E:333:PHE:H	1.82	0.45
1:F:106:ILE:HG23	1:F:107:TYR:CD1	2.51	0.45
1:F:200:PHE:CZ	1:F:390:CYS:HB3	2.52	0.45
1:G:125:CYS:O	1:G:125:CYS:SG	2.74	0.45
1:H:417:ALA:O	1:H:421:CYS:HB2	2.16	0.45
1:A:105:ARG:O	1:A:109:GLU:HG3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:364:PRO:HG2	1:C:369:CYS:SG	2.56	0.45
1:C:85:SER:CB	1:C:129:MET:HE2	2.46	0.45
1:D:359:ARG:HG2	1:D:360:PRO:HD2	1.98	0.45
1:A:245:LYS:HA	1:A:245:LYS:HD2	1.66	0.45
1:C:82:SER:HB2	1:C:189:SER:HG	1.75	0.45
1:E:6:ASP:HA	1:E:9:LYS:HG3	1.98	0.45
1:F:157:GLY:HA3	1:F:174:ALA:O	2.16	0.45
1:A:201:LYS:HD2	1:A:364:PRO:HA	1.98	0.45
1:B:245:LYS:HA	1:B:245:LYS:HD2	1.52	0.45
1:E:133:LEU:HD23	1:E:133:LEU:HA	1.80	0.45
1:E:352:ALA:O	1:E:356:VAL:HG23	2.17	0.45
1:G:279:ILE:HD12	1:G:296:ILE:HG22	1.98	0.45
1:C:86:HIS:CB	1:C:158:GLU:HB2	2.46	0.45
1:D:191:TYR:HD1	1:D:384:PRO:HB2	1.82	0.45
1:H:271:VAL:CG2	1:H:366:SER:HB3	2.45	0.45
1:B:231:VAL:HG12	1:B:232:ALA:O	2.17	0.45
1:C:124:ASN:HA	1:C:125:CYS:CB	2.43	0.45
1:C:38:GLU:CG	1:C:88:ASP:HB3	2.46	0.45
1:E:366:SER:N	1:E:367:PRO:CD	2.80	0.45
1:F:366:SER:N	1:F:367:PRO:CD	2.80	0.45
1:A:106:ILE:HG23	1:A:107:TYR:CD1	2.52	0.45
1:C:164:VAL:HG22	1:C:165:ASP:N	2.31	0.45
1:D:366:SER:N	1:D:367:PRO:CD	2.80	0.45
1:G:207:ALA:HB1	1:G:302:LEU:O	2.16	0.45
1:A:105:ARG:HG2	1:A:111:TRP:CZ3	2.51	0.44
1:A:349:LEU:HD22	1:A:386:LEU:HD11	1.99	0.44
1:B:107:TYR:HB2	1:B:108:HIS:CE1	2.52	0.44
1:B:162:GLU:HG3	1:H:185:HIS:HE1	1.81	0.44
1:F:114:GLU:OE1	1:F:115:GLY:N	2.50	0.44
1:H:243:MET:O	1:H:247:VAL:HG23	2.16	0.44
1:H:221:SER:OG	1:H:290:PHE:O	2.32	0.44
1:H:191:TYR:HD1	1:H:384:PRO:HB2	1.81	0.44
1:A:113:GLU:O	1:A:114:GLU:HB2	2.17	0.44
1:B:283:VAL:HA	1:B:284:PRO:HD3	1.84	0.44
1:C:216:VAL:HG12	1:C:243:MET:CE	2.47	0.44
1:C:392:GLY:O	1:C:398:ASN:HB3	2.17	0.44
1:D:212:LEU:HD12	1:D:300:ILE:HD13	1.92	0.44
1:G:6:ASP:O	1:G:9:LYS:HG3	2.17	0.44
1:A:106:ILE:HD11	1:A:399:THR:HB	2.00	0.44
1:E:76:GLY:O	1:E:147:LYS:HB3	2.17	0.44
1:F:352:ALA:O	1:F:356:VAL:HG23	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:77:GLY:O	1:F:424:THR:HG22	2.16	0.44
1:G:11:MET:SD	1:G:411:LYS:HG2	2.58	0.44
1:B:184:SER:O	1:H:332:LEU:HD12	2.17	0.44
1:B:114:GLU:C	1:B:116:ARG:N	2.67	0.44
1:B:123:VAL:HG12	1:B:394:ALA:HB2	1.99	0.44
1:B:274:VAL:HG12	1:B:275:ALA:N	2.33	0.44
1:C:201:LYS:HD2	1:C:364:PRO:HA	1.99	0.44
1:D:250:LEU:HD23	1:D:318:VAL:HG11	2.00	0.44
1:E:22:ILE:O	1:E:26:VAL:HG23	2.17	0.44
1:G:12:LYS:HD3	1:G:12:LYS:HA	1.64	0.44
1:G:164:VAL:HG22	1:G:165:ASP:N	2.32	0.44
1:H:211:PHE:HB2	1:H:332:LEU:HB3	1.98	0.44
1:A:366:SER:N	1:A:367:PRO:CD	2.81	0.44
1:B:86:HIS:CB	1:B:158:GLU:HB2	2.48	0.44
1:F:51:ARG:CG	1:F:51:ARG:NH1	2.81	0.44
1:A:224:THR:OG1	1:G:299:ASP:OD2	2.29	0.44
1:A:107:TYR:HB2	1:A:108:HIS:CE1	2.52	0.44
1:E:157:GLY:HA3	1:E:174:ALA:O	2.18	0.44
1:F:376:ASN:HB2	1:F:377:PRO:HD3	1.98	0.44
1:G:204:TRP:HB2	1:G:339:ALA:HB3	2.00	0.44
1:H:135:ALA:O	1:H:139:LEU:HG	2.17	0.44
1:H:260:ARG:HH22	1:H:314:GLU:CD	2.21	0.44
1:H:34:PRO:HG2	1:H:37:ARG:HD2	1.99	0.44
1:A:335:ARG:HB3	1:A:335:ARG:HE	1.63	0.44
1:A:364:PRO:HG2	1:A:369:CYS:SG	2.58	0.44
1:D:417:ALA:O	1:D:421:CYS:HB2	2.17	0.44
1:D:81:LEU:HD12	1:D:82:SER:H	1.82	0.44
1:H:124:ASN:HA	1:H:125:CYS:HB2	2.00	0.44
1:B:352:ALA:O	1:B:356:VAL:HG23	2.18	0.44
1:C:376:ASN:N	1:C:377:PRO:CD	2.81	0.44
1:D:394:ALA:O	1:D:397:GLY:HA3	2.18	0.44
1:D:34:PRO:HB3	1:D:92:ALA:HA	2.00	0.44
1:G:38:GLU:CG	1:G:88:ASP:HB3	2.47	0.44
1:H:86:HIS:HB3	1:H:158:GLU:HB2	1.99	0.44
1:H:88:ASP:OD1	1:H:88:ASP:N	2.49	0.44
1:A:101:ASP:OD1	1:A:101:ASP:N	2.51	0.44
1:A:211:PHE:HB2	1:A:332:LEU:HB3	1.99	0.44
1:B:34:PRO:HG2	1:B:37:ARG:HD2	2.00	0.44
1:C:123:VAL:HG12	1:C:394:ALA:HB2	2.00	0.44
1:C:195:ALA:O	1:C:196:GLU:HG2	2.18	0.44
1:D:28:LEU:HA	1:D:28:LEU:HD23	1.85	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:212:LEU:CD1	1:E:300:ILE:CD1	2.93	0.44
1:F:12:LYS:HD3	1:F:12:LYS:HA	1.66	0.44
1:G:110:ALA:HA	1:G:118:TYR:O	2.18	0.44
1:G:245:LYS:HD2	1:G:245:LYS:HA	1.69	0.44
1:A:240:ILE:CD1	1:A:291:PRO:CG	2.96	0.43
1:A:274:VAL:HG22	1:A:300:ILE:HG23	2.00	0.43
1:D:162:GLU:HG3	1:F:185:HIS:HE1	1.83	0.43
1:E:392:GLY:H	1:E:398:ASN:HD22	1.65	0.43
1:H:391:GLY:HA3	1:H:401:PHE:CE1	2.53	0.43
1:H:419:ASP:OD1	1:H:423:ARG:NE	2.46	0.43
1:A:210:VAL:HA	1:A:332:LEU:O	2.18	0.43
1:A:216:VAL:HG12	1:A:243:MET:HE1	2.00	0.43
1:C:91:MET:H	1:C:91:MET:HG2	1.68	0.43
1:E:240:ILE:CD1	1:E:291:PRO:CG	2.97	0.43
1:E:91:MET:HG2	1:E:91:MET:H	1.55	0.43
1:A:82:SER:HB2	1:A:189:SER:HG	1.80	0.43
1:C:224:THR:N	1:C:225:PRO:CD	2.81	0.43
1:C:63:ASP:N	1:C:63:ASP:OD1	2.40	0.43
1:D:107:TYR:HB2	1:D:108:HIS:CE1	2.53	0.43
1:E:243:MET:SD	1:E:296:ILE:HD13	2.58	0.43
1:B:212:LEU:HD22	1:B:328:VAL:HG12	1.94	0.43
1:B:26:VAL:O	1:B:30:SER:N	2.50	0.43
1:D:211:PHE:HB2	1:D:332:LEU:HB3	2.01	0.43
1:F:394:ALA:O	1:F:397:GLY:HA3	2.18	0.43
1:G:266:GLY:O	1:G:363:ARG:NH1	2.51	0.43
1:H:263:ARG:HB3	1:H:265:TYR:CE1	2.53	0.43
1:A:162:GLU:HG3	1:C:185:HIS:HE1	1.83	0.43
1:B:128:PRO:O	1:B:131:CYS:HB2	2.18	0.43
1:B:347:ASN:O	1:B:351:VAL:HG23	2.19	0.43
1:D:240:ILE:CD1	1:D:291:PRO:CG	2.96	0.43
1:D:352:ALA:O	1:D:356:VAL:HG23	2.19	0.43
1:C:319:VAL:HG11	1:C:326:ALA:HB3	2.00	0.43
1:G:179:ALA:O	1:G:183:ILE:HG13	2.19	0.43
1:G:212:LEU:CD2	1:G:328:VAL:HG12	2.48	0.43
1:H:155:VAL:HG11	1:H:179:ALA:HB2	2.01	0.43
1:H:183:ILE:O	1:H:185:HIS:N	2.52	0.43
1:B:192:ALA:O	1:B:385:SER:OG	2.33	0.43
1:C:274:VAL:HG22	1:C:300:ILE:HG23	2.00	0.43
1:C:347:ASN:O	1:C:351:VAL:HG23	2.18	0.43
1:C:366:SER:N	1:C:367:PRO:CD	2.82	0.43
1:E:212:LEU:HD12	1:E:300:ILE:HD13	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:300:ILE:HG21	1:E:311:VAL:HG11	2.01	0.43
1:G:376:ASN:N	1:G:377:PRO:CD	2.82	0.43
1:A:260:ARG:NH2	1:A:314:GLU:OE2	2.52	0.43
1:B:199:ASN:O	1:B:200:PHE:HB2	2.18	0.43
1:C:240:ILE:CD1	1:C:291:PRO:CG	2.97	0.43
1:F:347:ASN:O	1:F:351:VAL:HG23	2.19	0.43
1:G:258:GLU:O	1:G:262:THR:OG1	2.34	0.43
1:H:260:ARG:NH2	1:H:314:GLU:OE2	2.52	0.43
1:A:240:ILE:HD11	1:A:291:PRO:CG	2.48	0.43
1:B:307:ASN:ND2	1:B:309:LEU:HB2	2.34	0.43
1:B:34:PRO:HB3	1:B:92:ALA:HA	2.01	0.43
1:B:81:LEU:HD12	1:B:82:SER:N	2.34	0.43
1:C:86:HIS:H	1:C:86:HIS:CD2	2.36	0.43
1:E:18:ARG:O	1:E:22:ILE:HG13	2.19	0.43
1:F:183:ILE:HD11	1:F:378:TYR:CZ	2.54	0.43
1:F:343:GLU:HB3	1:F:344:PRO:CD	2.45	0.43
1:G:211:PHE:HB2	1:G:332:LEU:HB3	2.01	0.43
1:B:359:ARG:CG	1:B:360:PRO:HD2	2.43	0.42
1:F:164:VAL:HG22	1:F:165:ASP:N	2.34	0.42
1:B:44:TYR:CD1	1:B:44:TYR:C	2.92	0.42
1:C:85:SER:HB3	1:C:129:MET:HE2	2.01	0.42
1:E:7:VAL:HG22	1:E:418:MET:HE3	2.00	0.42
1:F:72:LEU:HB3	1:F:150:VAL:HB	2.00	0.42
1:F:216:VAL:HG12	1:F:243:MET:HE1	2.00	0.42
1:F:7:VAL:HG22	1:F:418:MET:CE	2.49	0.42
1:G:309:LEU:HD23	1:G:309:LEU:HA	1.78	0.42
1:G:352:ALA:O	1:G:356:VAL:HG23	2.18	0.42
1:B:238:ASN:HB3	1:B:241:VAL:HG23	2.02	0.42
1:D:116:ARG:HD3	1:D:400:TYR:CE2	2.53	0.42
1:D:191:TYR:OH	1:D:344:PRO:HG2	2.19	0.42
1:A:279:ILE:HD12	1:A:296:ILE:CG2	2.47	0.42
1:B:23:GLN:O	1:B:27:GLU:HG3	2.19	0.42
1:H:6:ASP:O	1:H:9:LYS:HG3	2.19	0.42
1:C:75:THR:HG23	1:C:146:LEU:O	2.19	0.42
1:E:207:ALA:HB1	1:E:302:LEU:O	2.20	0.42
1:G:263:ARG:HB3	1:G:265:TYR:HE1	1.85	0.42
1:H:12:LYS:HA	1:H:12:LYS:HD3	1.84	0.42
1:A:191:TYR:OH	1:A:344:PRO:HG2	2.20	0.42
1:B:300:ILE:HG21	1:B:311:VAL:HG11	2.01	0.42
1:B:38:GLU:CG	1:B:88:ASP:HB3	2.49	0.42
1:D:105:ARG:O	1:D:109:GLU:HG3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:12:LYS:HA	1:C:12:LYS:HD3	1.51	0.42
1:F:345:LEU:CD2	1:F:386:LEU:HG	2.49	0.42
1:H:114:GLU:C	1:H:116:ARG:N	2.70	0.42
1:A:81:LEU:HD12	1:A:82:SER:H	1.85	0.42
1:B:279:ILE:O	1:B:280:ARG:HB3	2.20	0.42
1:D:83:PHE:HE1	1:D:193:LEU:HD22	1.83	0.42
1:B:277:GLY:HA3	1:D:287:ILE:HB	2.02	0.42
1:E:110:ALA:HA	1:E:118:TYR:O	2.19	0.42
1:F:231:VAL:HG12	1:F:232:ALA:O	2.20	0.42
1:F:207:ALA:HB1	1:F:302:LEU:O	2.20	0.42
1:G:132:TRP:HZ3	1:G:152:LEU:HB3	1.83	0.42
1:H:265:TYR:OH	1:H:305:ASP:OD2	2.31	0.42
1:E:114:GLU:OE1	1:E:115:GLY:N	2.53	0.42
1:E:216:VAL:HG12	1:E:243:MET:HE1	2.02	0.42
1:E:231:VAL:CG1	1:E:232:ALA:N	2.82	0.42
1:F:88:ASP:N	1:F:88:ASP:OD1	2.51	0.42
1:G:105:ARG:O	1:G:109:GLU:HG3	2.20	0.42
1:G:204:TRP:HB2	1:G:339:ALA:CB	2.50	0.42
1:G:48:TRP:CE3	1:G:133:LEU:HD22	2.55	0.42
1:A:287:ILE:HB	1:G:277:GLY:HA3	2.02	0.42
1:B:12:LYS:HD3	1:B:12:LYS:HA	1.69	0.42
1:C:212:LEU:HD22	1:C:328:VAL:HG13	2.00	0.42
1:C:23:GLN:O	1:C:27:GLU:HG3	2.20	0.42
1:D:224:THR:O	1:D:227:VAL:HG12	2.20	0.42
2:E:501:6R6:O11	2:E:501:6R6:N13	2.52	0.42
1:F:216:VAL:HG12	1:F:243:MET:CE	2.49	0.42
1:G:374:ASP:O	1:G:377:PRO:HD2	2.19	0.42
1:H:23:GLN:O	1:H:27:GLU:HG3	2.20	0.42
1:A:34:PRO:HG2	1:A:37:ARG:HD2	2.02	0.41
1:A:45:VAL:O	1:A:49:MET:HB2	2.20	0.41
1:C:245:LYS:HD2	1:C:245:LYS:HA	1.64	0.41
1:C:256:ASN:O	1:C:260:ARG:HG3	2.20	0.41
1:C:7:VAL:HA	1:C:418:MET:HE1	2.01	0.41
1:D:12:LYS:HD3	1:D:12:LYS:HA	1.75	0.41
1:G:18:ARG:O	1:G:21:LEU:HB3	2.19	0.41
1:B:350:GLU:O	1:B:354:ARG:HG3	2.20	0.41
1:B:408:LYS:O	1:B:411:LYS:HB2	2.19	0.41
1:C:205:VAL:HG13	1:C:369:CYS:HB3	2.02	0.41
1:D:22:ILE:O	1:D:26:VAL:HG23	2.20	0.41
1:D:283:VAL:HA	1:D:284:PRO:HD3	1.88	0.41
1:D:46:TYR:C	1:D:46:TYR:CD1	2.94	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:243:MET:CE	1:E:296:ILE:HG23	2.38	0.41
1:F:86:HIS:ND1	1:F:158:GLU:HB3	2.34	0.41
1:G:81:LEU:HD12	1:G:82:SER:N	2.34	0.41
1:B:61:VAL:CG1	1:H:173:LEU:HD11	2.26	0.41
1:H:252:GLU:HG2	1:H:253:TRP:N	2.35	0.41
1:H:221:SER:HA	1:H:290:PHE:CE1	2.55	0.41
1:A:26:VAL:O	1:A:30:SER:OG	2.30	0.41
1:C:6:ASP:HA	1:C:9:LYS:HZ1	1.85	0.41
1:D:75:THR:HG23	1:D:146:LEU:O	2.21	0.41
1:H:212:LEU:CD1	1:H:300:ILE:CD1	2.94	0.41
1:A:199:ASN:O	1:A:200:PHE:HB2	2.21	0.41
1:A:283:VAL:HA	1:A:284:PRO:HD3	1.86	0.41
1:B:265:TYR:CZ	1:B:304:PRO:HD2	2.55	0.41
1:E:48:TRP:CZ3	1:E:133:LEU:HD22	2.55	0.41
1:H:349:LEU:HD22	1:H:386:LEU:HD11	2.02	0.41
1:A:195:ALA:O	1:A:196:GLU:HG2	2.20	0.41
1:B:366:SER:N	1:B:367:PRO:CD	2.84	0.41
1:B:91:MET:HG2	1:B:91:MET:H	1.70	0.41
1:C:208:GLY:N	1:C:302:LEU:O	2.52	0.41
1:D:240:ILE:CD1	1:D:291:PRO:HG2	2.50	0.41
1:E:231:VAL:HG12	1:E:232:ALA:O	2.20	0.41
1:F:260:ARG:HH22	1:F:314:GLU:CD	2.23	0.41
1:G:7:VAL:HG22	1:G:418:MET:CE	2.51	0.41
1:B:124:ASN:HB2	1:B:394:ALA:CB	2.51	0.41
1:D:51:ARG:CG	1:D:51:ARG:NH1	2.84	0.41
1:G:107:TYR:HB2	1:G:108:HIS:CE1	2.56	0.41
1:H:157:GLY:O	1:H:175:GLU:HA	2.20	0.41
1:B:129:MET:O	1:B:133:LEU:HB2	2.21	0.41
1:B:289:ARG:HD3	1:B:289:ARG:HH11	1.71	0.41
1:B:309:LEU:HA	1:B:309:LEU:HD23	1.81	0.41
1:B:124:ASN:HB2	1:B:394:ALA:HB2	2.02	0.41
1:C:107:TYR:HB2	1:C:108:HIS:CE1	2.56	0.41
1:E:302:LEU:HD21	1:E:311:VAL:HG21	2.02	0.41
1:F:206:GLU:HA	1:F:206:GLU:OE1	2.20	0.41
1:F:81:LEU:HD12	1:F:82:SER:N	2.35	0.41
1:H:240:ILE:HD11	1:H:291:PRO:CG	2.50	0.41
1:C:349:LEU:HD22	1:C:386:LEU:HD11	2.03	0.41
1:D:319:VAL:HG11	1:D:326:ALA:HB3	2.03	0.41
1:E:101:ASP:OD1	1:E:101:ASP:N	2.54	0.41
1:F:376:ASN:N	1:F:377:PRO:CD	2.84	0.41
1:F:46:TYR:C	1:F:46:TYR:CD1	2.94	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:332:LEU:HA	1:G:332:LEU:HD12	1.82	0.41
1:H:189:SER:O	1:H:383:ILE:HG12	2.20	0.41
1:A:114:GLU:OE1	1:A:115:GLY:N	2.54	0.41
1:A:125:CYS:O	1:A:125:CYS:SG	2.79	0.41
1:A:42:GLY:CA	1:A:87:LEU:HD13	2.50	0.41
1:B:224:THR:O	1:B:227:VAL:HG12	2.21	0.41
1:C:46:TYR:CD1	1:C:46:TYR:C	2.93	0.41
1:C:7:VAL:HG22	1:C:418:MET:HE3	2.02	0.41
1:D:350:GLU:O	1:D:354:ARG:HG3	2.21	0.41
1:D:91:MET:H	1:D:91:MET:HG2	1.63	0.41
1:E:195:ALA:O	1:E:196:GLU:HG2	2.21	0.41
1:E:47:GLU:O	1:E:51:ARG:HB2	2.21	0.41
1:H:107:TYR:HB2	1:H:108:HIS:CE1	2.56	0.41
1:A:175:GLU:OE1	1:A:209:LYS:NZ	2.44	0.41
1:A:91:MET:CE	1:A:120:TYR:CD1	3.03	0.41
1:B:210:VAL:HA	1:B:332:LEU:O	2.20	0.41
1:B:418:MET:O	1:B:422:ASN:ND2	2.53	0.41
1:F:191:TYR:OH	1:F:344:PRO:HG2	2.21	0.41
1:C:376:ASN:HB2	1:C:377:PRO:HD3	2.02	0.41
1:D:159:ILE:HG22	1:D:160:ASP:N	2.36	0.41
1:D:14:LEU:HD13	1:D:411:LYS:HG3	2.03	0.41
1:F:157:GLY:O	1:F:175:GLU:HA	2.21	0.41
1:H:212:LEU:HD22	1:H:328:VAL:HG12	2.03	0.41
1:A:158:GLU:HG2	1:A:374:ASP:OD2	2.20	0.40
1:B:211:PHE:O	1:B:330:PRO:HA	2.21	0.40
1:D:76:GLY:O	1:D:147:LYS:HB3	2.21	0.40
1:D:241:VAL:O	1:D:244:ALA:HB3	2.21	0.40
1:E:107:TYR:HB2	1:E:108:HIS:CE1	2.56	0.40
1:E:386:LEU:HD12	1:E:386:LEU:C	2.42	0.40
1:F:266:GLY:O	1:F:363:ARG:NH1	2.54	0.40
1:G:79:ALA:O	1:G:148:GLY:HA3	2.20	0.40
1:G:262:THR:CG2	1:G:271:VAL:HG22	2.51	0.40
1:H:14:LEU:CD2	1:H:138:ALA:HB2	2.51	0.40
1:A:88:ASP:OD2	1:A:159:ILE:HG13	2.20	0.40
1:C:160:ASP:HB2	1:H:288:TYR:HB3	2.04	0.40
1:C:170:HIS:ND1	1:C:170:HIS:C	2.74	0.40
1:C:343:GLU:HB3	1:C:344:PRO:HD3	2.04	0.40
1:G:366:SER:N	1:G:367:PRO:CD	2.83	0.40
1:H:196:GLU:O	1:H:389:GLY:HA3	2.20	0.40
1:B:124:ASN:HA	1:B:125:CYS:CB	2.52	0.40
1:A:329:LYS:NZ	1:C:57:GLU:OE1	2.37	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:211:PHE:O	1:D:330:PRO:HA	2.20	0.40
1:B:164:VAL:HG11	1:D:284:PRO:O	2.21	0.40
1:G:19:GLU:HG3	1:G:19:GLU:H	1.61	0.40
1:A:153:THR:HG22	1:A:155:VAL:CG1	2.50	0.40
1:A:164:VAL:HG22	1:A:165:ASP:N	2.36	0.40
1:E:347:ASN:O	1:E:351:VAL:HG23	2.22	0.40
1:G:133:LEU:HA	1:G:133:LEU:HD23	1.92	0.40
1:G:216:VAL:HG12	1:G:243:MET:HE1	2.03	0.40
1:H:101:ASP:N	1:H:101:ASP:OD1	2.54	0.40
1:H:164:VAL:HG22	1:H:165:ASP:N	2.37	0.40
1:H:332:LEU:HA	1:H:332:LEU:HD12	1.88	0.40
1:B:170:HIS:C	1:B:170:HIS:ND1	2.74	0.40
1:B:240:ILE:HD12	1:B:291:PRO:CD	2.51	0.40
1:B:28:LEU:HA	1:B:28:LEU:HD23	1.85	0.40
1:B:335:ARG:HE	1:B:335:ARG:HB3	1.77	0.40
1:F:227:VAL:HG13	1:F:227:VAL:O	2.22	0.40
1:E:251:GLU:CD	1:F:229:ARG:NH2	2.75	0.40
1:F:260:ARG:NH2	1:F:314:GLU:OE2	2.55	0.40
1:H:223:TYR:CG	1:H:225:PRO:HD2	2.57	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	418/425 (98%)	399 (96%)	17 (4%)	2 (0%)	34	77
1	B	418/425 (98%)	400 (96%)	17 (4%)	1 (0%)	52	88
1	C	418/425 (98%)	401 (96%)	16 (4%)	1 (0%)	52	88
1	D	418/425 (98%)	401 (96%)	16 (4%)	1 (0%)	52	88
1	E	418/425 (98%)	399 (96%)	18 (4%)	1 (0%)	52	88

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	418/425 (98%)	402 (96%)	15 (4%)	1 (0%)	52	88
1	G	418/425 (98%)	402 (96%)	14 (3%)	2 (0%)	34	77
1	H	418/425 (98%)	399 (96%)	17 (4%)	2 (0%)	34	77
All	All	3344/3400 (98%)	3203 (96%)	130 (4%)	11 (0%)	46	84

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	115	GLY
1	D	115	GLY
1	E	115	GLY
1	F	115	GLY
1	H	115	GLY
1	B	115	GLY
1	C	115	GLY
1	G	115	GLY
1	A	114	GLU
1	G	114	GLU
1	H	114	GLU

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	328/331 (99%)	313 (95%)	15 (5%)	33	73
1	B	328/331 (99%)	315 (96%)	13 (4%)	38	76
1	C	328/331 (99%)	313 (95%)	15 (5%)	33	73
1	D	328/331 (99%)	315 (96%)	13 (4%)	38	76
1	E	328/331 (99%)	316 (96%)	12 (4%)	41	78
1	F	328/331 (99%)	315 (96%)	13 (4%)	38	76
1	G	328/331 (99%)	315 (96%)	13 (4%)	38	76
1	H	328/331 (99%)	314 (96%)	14 (4%)	35	74

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	2624/2648 (99%)	2516 (96%)	108 (4%)	37 76

All (108) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ASP
1	A	17	MET
1	A	51	ARG
1	A	73	ARG
1	A	91	MET
1	A	94	GLU
1	A	95	ASP
1	A	101	ASP
1	A	109	GLU
1	A	160	ASP
1	A	162	GLU
1	A	216	VAL
1	A	359	ARG
1	A	362	GLU
1	A	402	LEU
1	B	17	MET
1	B	19	GLU
1	B	73	ARG
1	B	91	MET
1	B	95	ASP
1	B	109	GLU
1	B	159	ILE
1	B	160	ASP
1	B	162	GLU
1	B	216	VAL
1	B	362	GLU
1	B	387	THR
1	B	402	LEU
1	C	9	LYS
1	C	17	MET
1	C	18	ARG
1	C	51	ARG
1	C	67	ASN
1	C	73	ARG
1	C	91	MET
1	C	94	GLU
1	C	95	ASP

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Mol	Chain	Res	Type
1	C	109	GLU
1	C	160	ASP
1	C	162	GLU
1	C	216	VAL
1	C	362	GLU
1	C	402	LEU
1	D	6	ASP
1	D	9	LYS
1	D	17	MET
1	D	18	ARG
1	D	51	ARG
1	D	73	ARG
1	D	91	MET
1	D	95	ASP
1	D	109	GLU
1	D	160	ASP
1	D	162	GLU
1	D	216	VAL
1	D	402	LEU
1	E	6	ASP
1	E	17	MET
1	E	51	ARG
1	E	73	ARG
1	E	91	MET
1	E	95	ASP
1	E	109	GLU
1	E	160	ASP
1	E	162	GLU
1	E	216	VAL
1	E	362	GLU
1	E	402	LEU
1	F	6	ASP
1	F	17	MET
1	F	19	GLU
1	F	51	ARG
1	F	73	ARG
1	F	91	MET
1	F	95	ASP
1	F	109	GLU
1	F	160	ASP
1	F	162	GLU
1	F	216	VAL

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Mol	Chain	Res	Type
1	F	229	ARG
1	F	402	LEU
1	G	6	ASP
1	G	17	MET
1	G	19	GLU
1	G	51	ARG
1	G	73	ARG
1	G	91	MET
1	G	95	ASP
1	G	109	GLU
1	G	160	ASP
1	G	162	GLU
1	G	193	LEU
1	G	216	VAL
1	G	402	LEU
1	H	6	ASP
1	H	7	VAL
1	H	9	LYS
1	H	17	MET
1	H	51	ARG
1	H	73	ARG
1	H	91	MET
1	H	94	GLU
1	H	95	ASP
1	H	109	GLU
1	H	160	ASP
1	H	162	GLU
1	H	216	VAL
1	H	402	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	108	HIS
1	A	307	ASN
1	B	108	HIS
1	C	108	HIS
1	C	112	HIS
1	C	307	ASN
1	C	340	GLN
1	D	108	HIS
1	E	108	HIS

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Mol	Chain	Res	Type
1	E	307	ASN
1	F	108	HIS
1	F	340	GLN
1	G	108	HIS
1	G	307	ASN
1	H	108	HIS
1	H	307	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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	6R6	A	501	-	8,13,13	1.46	2 (25%)	11,18,18	1.17	1 (9%)
2	6R6	B	501	-	8,13,13	1.54	2 (25%)	11,18,18	1.67	2 (18%)
2	6R6	C	501	-	8,13,13	1.45	2 (25%)	11,18,18	1.45	3 (27%)
2	6R6	D	501	-	8,13,13	1.42	2 (25%)	11,18,18	2.19	4 (36%)
2	6R6	E	501	-	8,13,13	1.45	2 (25%)	11,18,18	1.29	2 (18%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	6R6	F	501	-	8,13,13	1.53	2 (25%)	11,18,18	2.14	3 (27%)
2	6R6	G	501	-	8,13,13	1.54	2 (25%)	11,18,18	1.41	1 (9%)
2	6R6	H	501	-	8,13,13	1.55	3 (37%)	11,18,18	1.09	1 (9%)

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	6R6	A	501	-	-	0/4/8/8	0/1/1/1
2	6R6	B	501	-	-	0/4/8/8	0/1/1/1
2	6R6	C	501	-	-	0/4/8/8	0/1/1/1
2	6R6	D	501	-	-	0/4/8/8	0/1/1/1
2	6R6	E	501	-	-	0/4/8/8	0/1/1/1
2	6R6	F	501	-	-	0/4/8/8	0/1/1/1
2	6R6	G	501	-	-	0/4/8/8	0/1/1/1
2	6R6	H	501	-	-	0/4/8/8	0/1/1/1

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	6R6	O08-N07	-2.90	1.17	1.22
2	B	501	6R6	O08-N07	-2.77	1.17	1.22
2	G	501	6R6	O08-N07	-2.76	1.17	1.22
2	E	501	6R6	O08-N07	-2.69	1.17	1.22
2	F	501	6R6	O08-N07	-2.65	1.17	1.22
2	C	501	6R6	O08-N07	-2.52	1.17	1.22
2	H	501	6R6	O08-N07	-2.46	1.17	1.22
2	D	501	6R6	O08-N07	-2.33	1.18	1.22
2	H	501	6R6	C04-C03	-2.23	1.38	1.41
2	A	501	6R6	C03-N13	2.21	1.44	1.37
2	E	501	6R6	C03-N13	2.26	1.45	1.37
2	H	501	6R6	C03-N13	2.32	1.45	1.37
2	G	501	6R6	C03-N13	2.36	1.45	1.37
2	F	501	6R6	C03-N13	2.41	1.45	1.37
2	D	501	6R6	C03-N13	2.48	1.45	1.37
2	B	501	6R6	C03-N13	2.55	1.46	1.37
2	C	501	6R6	C03-N13	2.58	1.46	1.37

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	6R6	C05-C06-N07	-2.27	116.75	118.74
2	C	501	6R6	C05-C04-C10	-2.12	116.92	120.16
2	D	501	6R6	C05-C04-C03	2.15	120.68	119.08
2	A	501	6R6	C05-C06-N07	2.19	120.67	118.74
2	E	501	6R6	C01-C06-N07	2.39	121.23	119.51
2	H	501	6R6	C05-C04-C03	2.49	120.93	119.08
2	C	501	6R6	C05-C04-C03	2.55	120.98	119.08
2	F	501	6R6	C05-C04-C03	2.56	120.99	119.08
2	E	501	6R6	C05-C04-C03	2.64	121.05	119.08
2	D	501	6R6	O08-N07-C06	2.79	120.68	118.67
2	B	501	6R6	C05-C04-C03	2.99	121.31	119.08
2	C	501	6R6	C01-C06-N07	3.14	121.77	119.51
2	B	501	6R6	C01-C06-N07	3.45	121.99	119.51
2	F	501	6R6	O08-N07-C06	3.53	121.21	118.67
2	G	501	6R6	C01-C06-N07	3.61	122.11	119.51
2	F	501	6R6	C01-C06-N07	4.96	123.07	119.51
2	D	501	6R6	C01-C06-N07	5.39	123.38	119.51

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	6R6	2	0
2	B	501	6R6	1	0
2	C	501	6R6	1	0
2	D	501	6R6	1	0
2	E	501	6R6	2	0
2	F	501	6R6	1	0
2	G	501	6R6	1	0
2	H	501	6R6	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	420/425 (98%)	-0.65	1 (0%) 95 94	9, 25, 59, 79	0
1	B	420/425 (98%)	-0.29	10 (2%) 62 50	12, 29, 81, 110	0
1	C	420/425 (98%)	-0.25	21 (5%) 32 21	9, 28, 86, 114	0
1	D	420/425 (98%)	-0.60	1 (0%) 95 94	11, 26, 66, 89	0
1	E	420/425 (98%)	-0.62	1 (0%) 95 94	10, 25, 59, 85	0
1	F	420/425 (98%)	-0.39	4 (0%) 84 77	11, 30, 67, 85	0
1	G	420/425 (98%)	-0.39	6 (1%) 78 67	10, 28, 70, 93	0
1	H	420/425 (98%)	-0.61	0 100 100	10, 26, 60, 82	0
All	All	3360/3400 (98%)	-0.48	44 (1%) 79 69	9, 27, 71, 114	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	113	GLU	4.5
1	B	116	ARG	4.1
1	C	114	GLU	4.1
1	B	114	GLU	4.1
1	C	116	ARG	3.7
1	B	113	GLU	3.5
1	B	115	GLY	3.4
1	C	118	TYR	3.4
1	B	399	THR	3.3
1	C	398	ASN	3.3
1	B	118	TYR	3.3
1	C	400	TYR	3.2
1	C	103	ASN	3.2
1	B	106	ILE	3.0
1	C	397	GLY	3.0
1	C	107	TYR	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	22	ILE	2.8
1	C	106	ILE	2.7
1	C	402	LEU	2.7
1	G	113	GLU	2.7
1	B	400	TYR	2.7
1	F	399	THR	2.6
1	C	104	ASP	2.6
1	C	108	HIS	2.6
1	C	394	ALA	2.5
1	B	401	PHE	2.5
1	G	115	GLY	2.5
1	C	396	GLY	2.4
1	F	114	GLU	2.4
1	G	398	ASN	2.4
1	G	114	GLU	2.4
1	C	109	GLU	2.3
1	C	401	PHE	2.3
1	C	399	THR	2.3
1	C	393	GLY	2.3
1	C	101	ASP	2.3
1	G	16	GLY	2.3
1	F	113	GLU	2.2
1	B	404	ASP	2.2
1	D	106	ILE	2.1
1	G	404	ASP	2.0
1	F	8	ALA	2.0
1	E	104	ASP	2.0
1	A	113	GLU	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 carbohydrates 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. LLDF column lists the quality of electron

density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	6R6	F	501	13/13	0.96	0.26	3.19	37,44,47,49	0
2	6R6	D	501	13/13	0.93	0.25	2.50	35,48,57,60	0
2	6R6	G	501	13/13	0.95	0.25	2.32	34,44,52,53	0
2	6R6	A	501	13/13	0.95	0.22	2.14	28,34,42,43	0
2	6R6	E	501	13/13	0.93	0.23	1.96	30,38,44,45	0
2	6R6	C	501	13/13	0.86	0.38	1.47	49,61,74,75	0
2	6R6	H	501	13/13	0.96	0.22	1.43	35,43,55,59	0
2	6R6	B	501	13/13	0.90	0.25	0.80	43,47,64,64	0

6.5 Other polymers ⓘ

There are no such residues in this entry.