



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:15 AM GMT

PDB ID : 2C8N
Title : THE STRUCTURE OF A FAMILY 51 ARABINOFURANOSIDASE, ARAF51, FROM CLOSTRIDIUM THERMOCELLUM IN COMPLEX WITH 1,3-LINKED ARABINOSIDE OF XYLOBIOSE.
Authors : Taylor, E.J.; Smith, N.L.; Turkenburg, J.P.; D'Souza, S.; Gilbert, H.J.; Davies, G.J.
Deposited on : 2005-12-06
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
<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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
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) : trunk26865

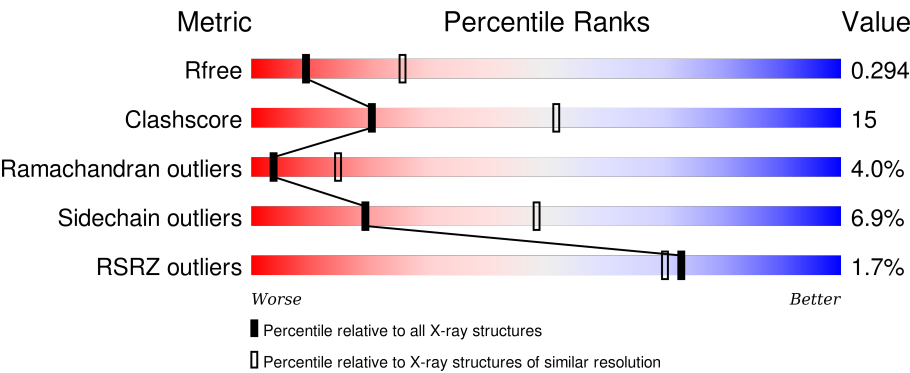
1 Overall quality at a glance ⓘ

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}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (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 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	513	<div><div>2%</div><div>77%</div><div>16%</div><div>• • •</div></div>
1	B	513	<div><div>%</div><div>69%</div><div>21%</div><div>5%</div><div>• • •</div></div>
1	C	513	<div><div>2%</div><div>71%</div><div>22%</div><div>• • •</div></div>
1	D	513	<div><div>%</div><div>71%</div><div>21%</div><div>5%</div><div>• • •</div></div>
1	E	513	<div><div>2%</div><div>73%</div><div>18%</div><div>6%</div><div>•</div></div>

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Mol	Chain	Length	Quality of chain
1	F	513	

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
3	EDO	A	1503	-	-	-	X
3	EDO	B	1503	-	-	-	X
3	EDO	C	1503	-	-	-	X
3	EDO	E	1503	-	-	-	X
3	EDO	F	1503	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 24200 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 ALPHA-L-ARABINOFURANOSIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	501	Total	C	N	O	S	0	0	0
			4017	2552	683	760	22			
1	B	500	Total	C	N	O	S	0	0	0
			4013	2545	686	760	22			
1	C	499	Total	C	N	O	S	0	1	0
			3989	2533	681	753	22			
1	D	498	Total	C	N	O	S	0	0	0
			3933	2487	672	753	21			
1	E	497	Total	C	N	O	S	0	0	0
			3927	2496	670	741	20			
1	F	500	Total	C	N	O	S	0	0	0
			3966	2518	679	748	21			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	173	ALA	GLU	ENGINEERED MUTATION	UNP Q4CJG5
B	173	ALA	GLU	ENGINEERED MUTATION	UNP Q4CJG5
C	173	ALA	GLU	ENGINEERED MUTATION	UNP Q4CJG5
D	173	ALA	GLU	ENGINEERED MUTATION	UNP Q4CJG5
E	173	ALA	GLU	ENGINEERED MUTATION	UNP Q4CJG5
F	173	ALA	GLU	ENGINEERED MUTATION	UNP Q4CJG5

- Molecule 2 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	2	Total	C	O	0	0
			18	10	8		
2	B	2	Total	C	O	0	0
			18	10	8		
2	C	2	Total	C	O	0	0
			18	10	8		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	2	Total	C	O	0	0
			18	10	8		
2	E	2	Total	C	O	0	0
			18	10	8		
2	F	2	Total	C	O	0	0
			18	10	8		

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	E	1	Total	C	O	0	0
			4	2	2		
3	E	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	F	1	Total	C	O	0	0
			4	2	2		

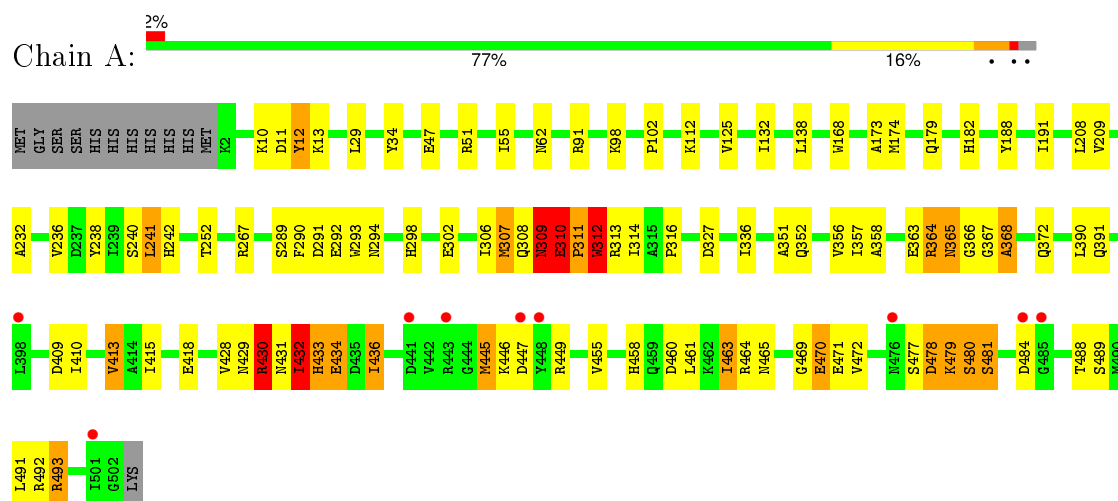
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	35	Total	O	0	0
			35	35		
4	B	37	Total	O	0	0
			37	37		
4	C	36	Total	O	0	0
			36	36		
4	D	31	Total	O	0	0
			31	31		
4	E	40	Total	O	0	0
			40	40		
4	F	32	Total	O	0	0
			32	32		

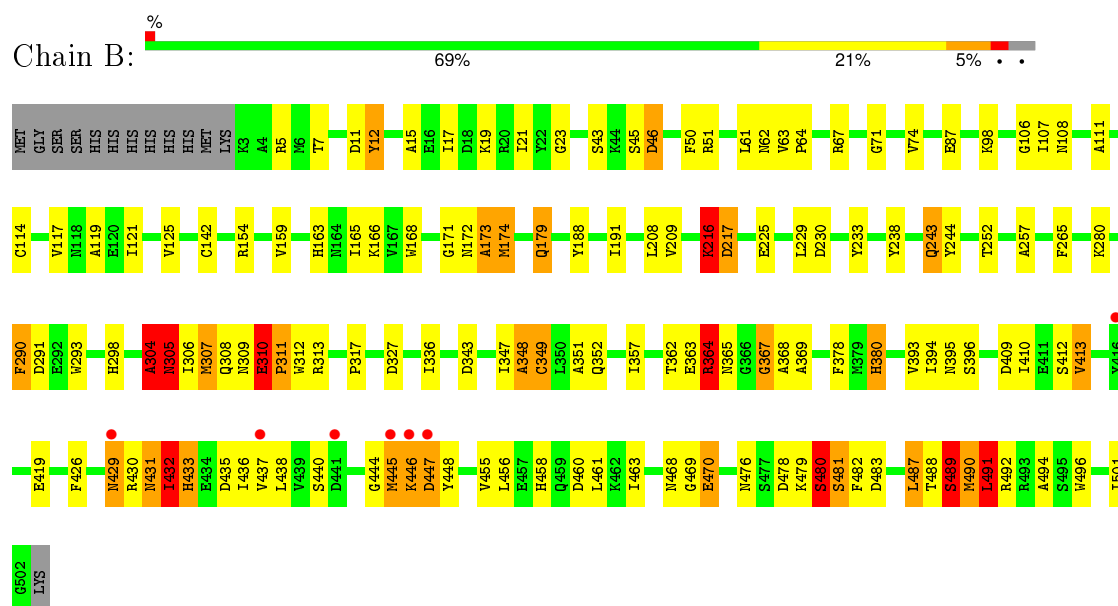
3 Residue-property plots

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: ALPHA-L-ARABINOFURANOSIDASE

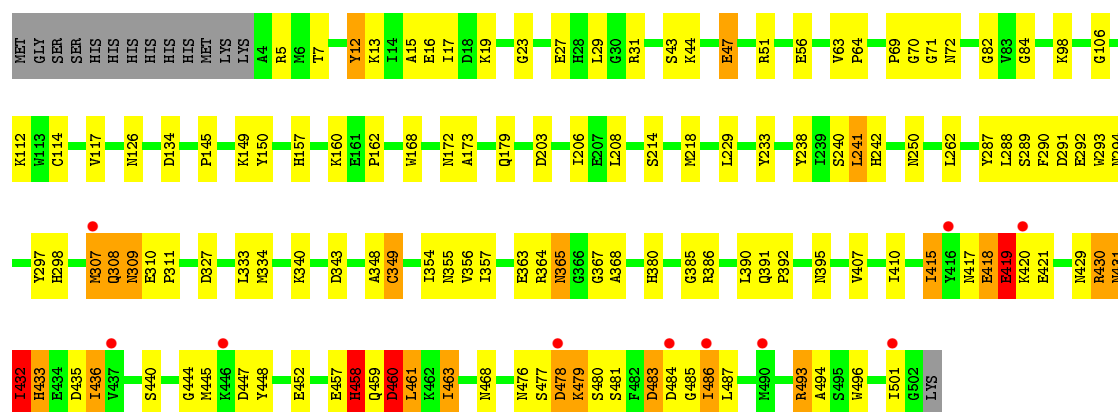


• Molecule 1: ALPHA-L-ARABINOFURANOSIDASE

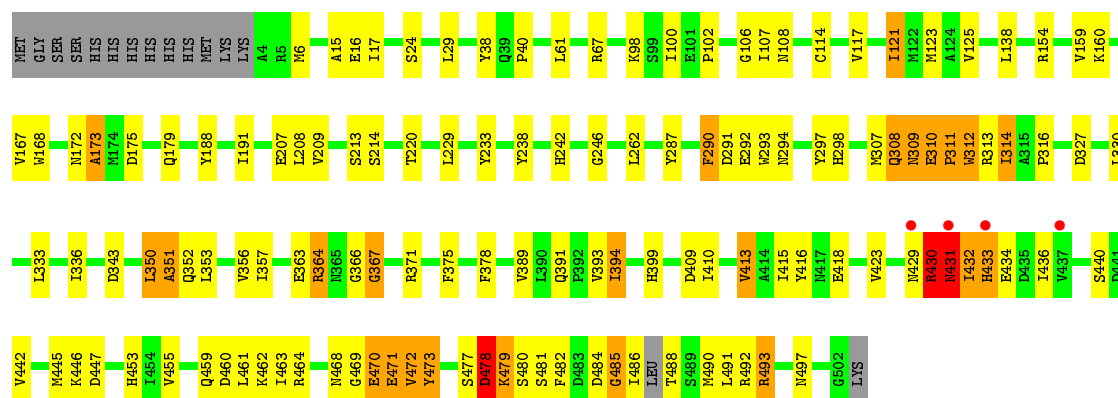


• Molecule 1: ALPHA-L-ARABINOFURANOSIDASE

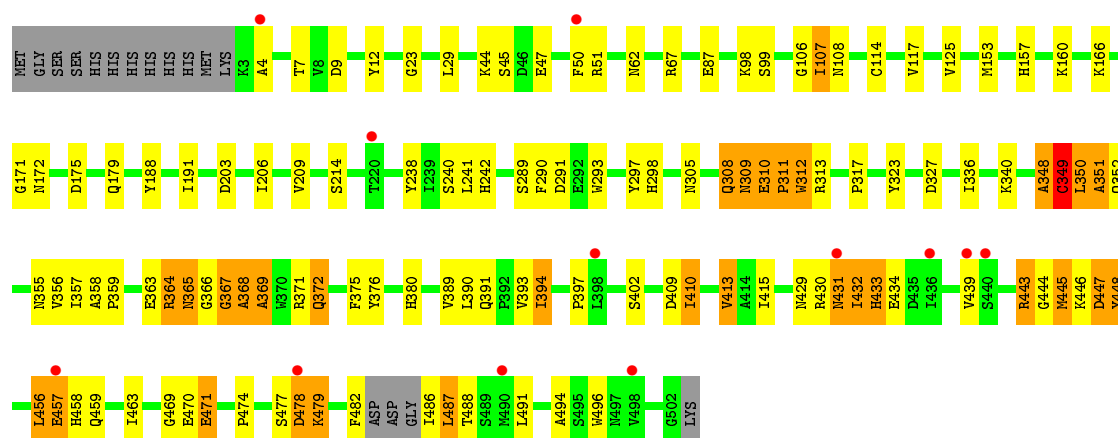
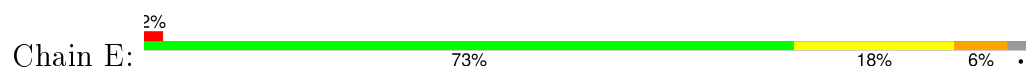




• Molecule 1: ALPHA-L-ARABINOFURANOSIDASE



• Molecule 1: ALPHA-L-ARABINOFURANOSIDASE



• Molecule 1: ALPHA-L-ARABINOFURANOSIDASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	173.45Å 173.45Å 272.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	145.86 – 2.90 63.35 – 2.43	Depositor EDS
% Data completeness (in resolution range)	98.8 (145.86-2.90) 92.4 (63.35-2.43)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.25 (at 2.42Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.201 , 0.255 0.250 , 0.294	Depositor DCC
R_{free} test set	4555 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	54.5	Xtriage
Anisotropy	0.066	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 53.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	1 of 144362 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	24200	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.05% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: AHR, XYS, EDO

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	0.38	0/4110	0.61	1/5566 (0.0%)
1	B	0.37	0/4105	0.60	0/5558
1	C	0.38	0/4082	0.60	0/5531
1	D	0.37	0/4021	0.59	0/5447
1	E	0.39	1/4019 (0.0%)	0.61	0/5447
1	F	0.37	0/4058	0.56	0/5496
All	All	0.38	1/24395 (0.0%)	0.60	1/33045 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	5
1	B	0	6
1	C	0	2
1	D	0	4
1	E	0	3
All	All	0	20

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	402	SER	CB-OG	5.30	1.49	1.42

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	310	GLU	C-N-CD	-9.15	100.47	120.60

There are no chirality outliers.

All (20) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	309	ASN	Peptide
1	A	310	GLU	Peptide
1	A	312	TRP	Peptide
1	A	432	ILE	Peptide
1	A	433	HIS	Peptide
1	B	216	LYS	Peptide
1	B	304	ALA	Peptide
1	B	347	ILE	Peptide
1	B	348	ALA	Peptide
1	B	45	SER	Peptide
1	B	491	LEU	Peptide
1	C	419	GLU	Peptide
1	C	458	HIS	Peptide
1	D	312	TRP	Peptide
1	D	431	ASN	Peptide
1	D	471	GLU	Peptide
1	D	485	GLY	Peptide
1	E	308	GLN	Peptide
1	E	348	ALA	Peptide
1	E	459	GLN	Peptide

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	4017	0	3920	117	0
1	B	4013	0	3908	158	0
1	C	3989	0	3870	107	0
1	D	3933	0	3767	113	0
1	E	3927	0	3786	109	0
1	F	3966	0	3845	90	0
2	A	18	0	15	0	0
2	B	18	0	15	1	0
2	C	18	0	15	0	0
2	D	18	0	15	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	18	0	15	0	0
2	F	18	0	15	0	0
3	A	8	0	12	0	0
3	B	4	0	6	0	0
3	C	8	0	12	0	0
3	D	4	0	6	0	0
3	E	8	0	12	0	0
3	F	4	0	6	0	0
4	A	35	0	0	1	0
4	B	37	0	0	1	0
4	C	36	0	0	0	0
4	D	31	0	0	0	0
4	E	40	0	0	1	0
4	F	32	0	0	1	0
All	All	24200	0	23240	691	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 (691) 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:431:ASN:CB	1:B:432:ILE:HG22	1.11	1.56
1:D:431:ASN:HB3	1:D:432:ILE:C	1.38	1.40
1:B:431:ASN:CB	1:B:432:ILE:CG2	2.05	1.32
1:A:310:GLU:HB3	1:A:311:PRO:CB	1.73	1.18
1:E:478:ASP:HA	1:E:479:LYS:HB2	1.22	1.18
1:B:431:ASN:HA	1:B:432:ILE:HB	1.18	1.17
1:F:469:GLY:HA2	1:F:470:GLU:HB2	1.17	1.16
1:A:310:GLU:HB3	1:A:311:PRO:HB3	1.20	1.12
1:C:410:ILE:HD11	1:C:436:ILE:HG13	1.32	1.11
1:C:478:ASP:HB3	1:C:479:LYS:HA	1.23	1.11
1:B:480:SER:HA	1:B:481:SER:HB2	1.33	1.10
1:A:469:GLY:HA2	1:A:470:GLU:HB2	1.20	1.10
1:D:480:SER:HB3	1:D:481:SER:HA	1.24	1.09
1:A:433:HIS:H	1:A:434:GLU:HB3	1.15	1.09
1:F:367:GLY:HA3	1:F:368:ALA:HB2	1.35	1.08
1:E:367:GLY:HA2	1:E:368:ALA:CB	1.81	1.08
1:B:307:MET:HG3	1:B:308:GLN:H	1.11	1.07
1:E:447:ASP:HB2	1:E:448:TYR:HA	1.08	1.07
1:E:447:ASP:CB	1:E:448:TYR:HA	1.85	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:348:ALA:HA	1:E:349:CYS:HB2	1.32	1.06
1:B:431:ASN:HB2	1:B:432:ILE:CG2	1.74	1.06
1:D:469:GLY:HA2	1:D:470:GLU:HB3	1.36	1.06
1:A:367:GLY:HA3	1:A:368:ALA:CB	1.83	1.06
1:B:429:ASN:CG	1:B:491:LEU:HD12	1.76	1.05
1:A:480:SER:HA	1:A:481:SER:HB2	1.08	1.05
1:A:310:GLU:H	1:A:311:PRO:HD3	1.21	1.05
1:F:429:ASN:HA	1:F:430:ARG:O	1.57	1.03
1:C:420:LYS:N	1:C:421:GLU:HA	1.70	1.03
1:E:445:MET:HA	1:E:446:LYS:HB3	1.36	1.03
1:D:431:ASN:HB3	1:D:432:ILE:O	1.58	1.02
1:B:431:ASN:HB3	1:B:432:ILE:CG2	1.76	1.01
1:A:433:HIS:H	1:A:434:GLU:CB	1.74	1.01
1:D:350:LEU:HA	1:D:351:ALA:HB3	1.42	1.00
1:B:431:ASN:HB3	1:B:432:ILE:HG22	1.00	0.99
1:D:430:ARG:HA	1:D:431:ASN:O	1.61	0.99
1:B:304:ALA:HB3	1:B:305:ASN:HB2	1.41	0.99
1:D:311:PRO:HG2	1:D:312:TRP:NE1	1.77	0.99
1:B:431:ASN:HA	1:B:432:ILE:CB	1.91	0.99
1:F:367:GLY:CA	1:F:368:ALA:HB2	1.93	0.99
1:A:410:ILE:HD11	1:A:436:ILE:HD11	1.45	0.99
1:A:367:GLY:HA3	1:A:368:ALA:HB2	0.99	0.99
1:A:367:GLY:CA	1:A:368:ALA:HB2	1.92	0.98
1:A:310:GLU:N	1:A:311:PRO:HD3	1.74	0.98
1:F:469:GLY:CA	1:F:470:GLU:HB2	1.93	0.98
1:B:431:ASN:HB2	1:B:432:ILE:HG22	0.97	0.96
1:C:348:ALA:HA	1:C:349:CYS:HB2	1.47	0.96
1:B:432:ILE:HG23	1:B:432:ILE:O	1.64	0.95
1:C:431:ASN:HA	1:C:432:ILE:O	1.65	0.95
1:C:363:GLU:HB2	1:C:368:ALA:HB3	1.47	0.95
1:B:348:ALA:CA	1:B:349:CYS:HB2	1.98	0.94
1:E:367:GLY:HA2	1:E:368:ALA:HB2	1.47	0.94
1:A:469:GLY:CA	1:A:470:GLU:HB2	1.97	0.94
1:D:480:SER:CB	1:D:481:SER:HA	1.97	0.94
1:E:368:ALA:HA	1:E:369:ALA:HB2	1.50	0.93
1:E:350:LEU:HA	1:E:351:ALA:HB3	1.51	0.93
1:B:431:ASN:CA	1:B:432:ILE:HB	1.99	0.92
1:A:364:ARG:C	1:A:366:GLY:HA3	1.91	0.91
1:D:471:GLU:O	1:D:472:VAL:HG23	1.69	0.91
1:A:363:GLU:H	1:A:368:ALA:HB1	1.36	0.90
1:C:363:GLU:HB2	1:C:368:ALA:CB	2.01	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:431:ASN:CB	1:D:432:ILE:C	2.34	0.90
1:D:351:ALA:H	1:D:352:GLN:HA	1.37	0.90
1:E:447:ASP:HB2	1:E:448:TYR:CA	2.00	0.89
1:B:348:ALA:CB	1:B:349:CYS:HB2	2.03	0.89
1:E:351:ALA:H	1:E:352:GLN:HA	1.38	0.88
1:A:477:SER:HB3	1:A:478:ASP:HA	1.53	0.88
1:B:431:ASN:CA	1:B:432:ILE:CB	2.51	0.88
1:A:310:GLU:CB	1:A:311:PRO:HB3	2.03	0.88
1:A:480:SER:HA	1:A:481:SER:CB	1.96	0.88
1:D:312:TRP:N	1:D:313:ARG:HB2	1.89	0.87
1:B:469:GLY:HA2	1:B:470:GLU:CB	2.04	0.87
1:A:477:SER:HB3	1:A:478:ASP:CA	2.05	0.87
1:C:419:GLU:HB3	1:C:420:LYS:HB3	1.54	0.87
1:D:311:PRO:HG2	1:D:312:TRP:CD1	2.09	0.87
1:B:307:MET:CG	1:B:308:GLN:H	1.87	0.87
1:E:478:ASP:CA	1:E:479:LYS:HB2	2.04	0.86
1:A:309:ASN:CB	1:A:310:GLU:HA	2.05	0.86
1:B:491:LEU:HB3	1:B:492:ARG:O	1.73	0.86
1:F:469:GLY:HA2	1:F:470:GLU:CB	2.04	0.85
1:E:291:ASP:HA	1:E:349:CYS:HB3	1.58	0.85
1:B:348:ALA:HB1	1:B:349:CYS:HB2	1.58	0.85
1:E:310:GLU:H	1:E:311:PRO:HA	1.41	0.85
1:E:457:GLU:HB3	1:E:458:HIS:HB2	1.57	0.85
1:C:460:ASP:HA	1:C:461:LEU:CB	2.05	0.85
1:D:469:GLY:HA2	1:D:470:GLU:CB	2.07	0.85
1:E:367:GLY:HA2	1:E:368:ALA:HB3	1.58	0.85
1:B:429:ASN:CG	1:B:491:LEU:CD1	2.45	0.84
1:B:216:LYS:N	1:B:217:ASP:HB2	1.92	0.83
1:B:304:ALA:CB	1:B:305:ASN:HB2	2.08	0.83
1:E:457:GLU:HB3	1:E:458:HIS:CA	2.08	0.83
1:B:429:ASN:OD1	1:B:491:LEU:HD12	1.78	0.83
1:B:348:ALA:HA	1:B:349:CYS:HB2	1.61	0.83
1:A:469:GLY:HA2	1:A:470:GLU:CB	2.07	0.83
1:E:350:LEU:HA	1:E:351:ALA:CB	2.09	0.83
1:E:308:GLN:N	1:E:309:ASN:O	2.11	0.83
1:D:350:LEU:HA	1:D:351:ALA:CB	2.08	0.83
1:C:291:ASP:HA	1:C:349:CYS:HB3	1.61	0.82
1:A:309:ASN:HB2	1:A:310:GLU:HA	1.60	0.82
1:A:465:ASN:HD21	1:A:472:VAL:H	1.28	0.81
1:A:433:HIS:N	1:A:434:GLU:HB3	1.96	0.81
1:E:348:ALA:CA	1:E:349:CYS:HB2	2.11	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:410:ILE:HD11	1:C:436:ILE:CG1	2.10	0.81
1:A:312:TRP:N	1:A:313:ARG:HB2	1.94	0.81
1:E:308:GLN:CB	1:E:309:ASN:HB2	2.10	0.81
1:B:367:GLY:N	1:B:368:ALA:HB2	1.96	0.80
1:C:478:ASP:CB	1:C:479:LYS:HA	2.04	0.80
1:C:420:LYS:H	1:C:421:GLU:HA	1.44	0.80
1:B:431:ASN:CA	1:B:432:ILE:HG22	2.11	0.79
1:C:410:ILE:CD1	1:C:436:ILE:HG13	2.12	0.79
1:A:363:GLU:H	1:A:368:ALA:CB	1.95	0.79
1:E:457:GLU:HB3	1:E:458:HIS:CB	2.11	0.79
1:C:432:ILE:HG22	1:C:494:ALA:HB2	1.65	0.79
1:B:482:PHE:HB2	1:B:487:LEU:HD12	1.63	0.79
1:D:366:GLY:HA3	1:D:367:GLY:O	1.82	0.79
1:B:348:ALA:HB1	1:B:349:CYS:CB	2.13	0.78
1:E:470:GLU:CB	1:E:471:GLU:HB2	2.14	0.78
1:E:363:GLU:O	1:E:364:ARG:HB2	1.84	0.78
1:B:469:GLY:HA2	1:B:470:GLU:HB3	1.65	0.78
1:B:409:ASP:O	1:B:430:ARG:HB3	1.82	0.78
1:F:188:TYR:HA	1:F:191:ILE:HG22	1.64	0.78
1:B:367:GLY:CA	1:B:368:ALA:HB2	2.15	0.77
1:D:480:SER:HB3	1:D:481:SER:CA	2.11	0.77
1:B:432:ILE:CG2	1:B:432:ILE:O	2.31	0.76
1:D:429:ASN:HA	1:D:430:ARG:O	1.85	0.76
1:F:478:ASP:HA	1:F:479:LYS:HG2	1.67	0.76
1:B:431:ASN:HB2	1:B:432:ILE:C	2.06	0.76
1:F:310:GLU:CB	1:F:311:PRO:HA	2.16	0.76
1:E:366:GLY:HA2	1:E:367:GLY:O	1.85	0.76
1:C:291:ASP:HA	1:C:349:CYS:CB	2.16	0.76
1:D:472:VAL:O	1:D:473:TYR:HB3	1.86	0.76
1:F:380:HIS:HD2	1:F:496:TRP:HE1	1.32	0.75
1:E:367:GLY:CA	1:E:368:ALA:CB	2.64	0.75
1:B:307:MET:HG3	1:B:308:GLN:N	1.96	0.74
1:B:216:LYS:H	1:B:217:ASP:HB2	1.51	0.74
1:E:445:MET:HA	1:E:446:LYS:CB	2.17	0.74
1:A:310:GLU:HB3	1:A:311:PRO:CG	2.17	0.74
1:B:291:ASP:HA	1:B:349:CYS:CB	2.18	0.74
1:B:458:HIS:HD2	1:B:460:ASP:H	1.35	0.73
1:D:431:ASN:HB3	1:D:432:ILE:CA	2.18	0.73
1:C:432:ILE:HG13	1:C:432:ILE:O	1.86	0.73
1:B:431:ASN:CA	1:B:432:ILE:CG2	2.66	0.73
1:C:478:ASP:HB3	1:C:479:LYS:CA	2.11	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:420:LYS:N	1:C:421:GLU:CA	2.50	0.73
1:F:367:GLY:CA	1:F:368:ALA:CB	2.65	0.73
1:A:480:SER:CA	1:A:481:SER:HB2	2.02	0.73
1:C:418:GLU:H	1:C:419:GLU:C	1.91	0.73
1:B:310:GLU:HG2	1:B:313:ARG:HD2	1.70	0.72
1:A:410:ILE:HD13	1:A:429:ASN:HB2	1.71	0.72
1:B:491:LEU:CB	1:B:492:ARG:CA	2.66	0.72
1:E:431:ASN:HD22	1:E:431:ASN:H	1.38	0.72
1:A:433:HIS:N	1:A:434:GLU:CB	2.51	0.72
1:E:351:ALA:N	1:E:352:GLN:HA	2.00	0.72
1:C:460:ASP:HA	1:C:461:LEU:HB2	1.70	0.72
1:D:309:ASN:N	1:D:310:GLU:HB2	2.05	0.72
1:B:478:ASP:HA	1:B:479:LYS:HB2	1.71	0.71
1:E:291:ASP:HA	1:E:349:CYS:CB	2.20	0.70
1:B:310:GLU:HB3	1:B:311:PRO:HA	1.72	0.70
1:E:340:LYS:HG3	1:E:413:VAL:HG11	1.73	0.70
1:B:46:ASP:HB3	1:B:50:PHE:O	1.90	0.70
1:B:431:ASN:H	1:B:431:ASN:ND2	1.89	0.70
1:B:446:LYS:HD3	1:B:446:LYS:H	1.57	0.70
1:D:431:ASN:CB	1:D:432:ILE:O	2.39	0.70
1:E:447:ASP:CB	1:E:448:TYR:CA	2.63	0.70
1:C:431:ASN:HA	1:C:432:ILE:C	2.13	0.70
1:A:312:TRP:CA	1:A:313:ARG:HB2	2.22	0.69
1:E:363:GLU:H	1:E:368:ALA:H	1.38	0.69
1:D:312:TRP:H	1:D:313:ARG:HB2	1.54	0.69
1:A:309:ASN:HB2	1:A:310:GLU:CA	2.22	0.69
1:F:51:ARG:HH22	1:F:368:ALA:HB3	1.56	0.69
1:A:460:ASP:HB3	1:A:463:ILE:HB	1.74	0.69
1:B:304:ALA:CA	1:B:305:ASN:HB2	2.22	0.69
1:B:291:ASP:HA	1:B:349:CYS:HB3	1.74	0.69
1:C:363:GLU:O	1:C:364:ARG:HB3	1.93	0.69
1:C:435:ASP:OD1	1:C:493:ARG:HG3	1.93	0.69
1:B:436:ILE:O	1:B:491:LEU:HG	1.93	0.68
1:C:15:ALA:HB2	1:C:343:ASP:HB3	1.74	0.68
1:A:460:ASP:O	1:A:461:LEU:HB2	1.94	0.68
1:D:351:ALA:N	1:D:352:GLN:HA	1.99	0.68
1:C:333:LEU:HD22	1:C:430:ARG:HH11	1.59	0.67
1:E:457:GLU:HB3	1:E:458:HIS:HA	1.75	0.67
1:B:216:LYS:CE	1:B:216:LYS:O	2.41	0.67
1:B:23:GLY:HA2	1:B:63:VAL:HG13	1.76	0.67
1:B:310:GLU:CB	1:B:311:PRO:HA	2.24	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:432:ILE:O	1:B:433:HIS:ND1	2.29	0.66
1:B:491:LEU:HB2	1:B:492:ARG:HA	1.78	0.66
1:E:308:GLN:CB	1:E:309:ASN:CB	2.73	0.66
1:A:298:HIS:HE1	1:A:327:ASP:OD2	1.77	0.66
1:A:312:TRP:HA	1:A:313:ARG:HB2	1.77	0.66
1:A:477:SER:HB3	1:A:478:ASP:CB	2.26	0.66
1:C:418:GLU:N	1:C:419:GLU:C	2.49	0.66
1:C:460:ASP:HA	1:C:461:LEU:HB3	1.78	0.66
1:D:15:ALA:HB2	1:D:343:ASP:HB3	1.77	0.66
1:B:480:SER:HA	1:B:481:SER:CB	2.14	0.66
1:C:348:ALA:CA	1:C:349:CYS:HB2	2.24	0.66
1:D:478:ASP:HA	1:D:479:LYS:CB	2.26	0.66
1:D:29:LEU:HD13	1:D:356:VAL:HG11	1.78	0.66
1:C:172:ASN:HD22	1:C:179:GLN:HE22	1.40	0.66
1:F:363:GLU:OE1	1:F:363:GLU:HA	1.96	0.65
1:F:478:ASP:HA	1:F:479:LYS:CG	2.26	0.65
1:A:310:GLU:HB3	1:A:311:PRO:CD	2.25	0.65
1:A:364:ARG:O	1:A:366:GLY:HA3	1.96	0.65
1:F:310:GLU:HB3	1:F:311:PRO:HA	1.78	0.65
1:F:458:HIS:HD2	1:F:460:ASP:H	1.41	0.65
1:A:410:ILE:CD1	1:A:436:ILE:HD11	2.24	0.64
1:D:307:MET:C	1:D:308:GLN:HG3	2.18	0.64
1:A:363:GLU:O	1:A:364:ARG:HG2	1.97	0.64
1:B:491:LEU:CB	1:B:492:ARG:HA	2.28	0.64
1:B:230:ASP:O	1:B:280:LYS:NZ	2.28	0.64
1:C:477:SER:O	1:C:478:ASP:HB2	1.97	0.64
1:C:31:ARG:NH1	1:C:364:ARG:HB2	2.13	0.64
1:A:365:ASN:N	1:A:366:GLY:HA3	2.10	0.64
1:B:480:SER:CA	1:B:481:SER:HB2	2.20	0.63
1:F:380:HIS:CD2	1:F:496:TRP:HE1	2.14	0.63
1:D:482:PHE:HA	1:D:486:ILE:O	1.98	0.63
1:B:310:GLU:HB3	1:B:311:PRO:CA	2.29	0.63
1:B:431:ASN:CB	1:B:432:ILE:CB	2.74	0.63
1:E:445:MET:CA	1:E:446:LYS:HB3	2.19	0.63
1:F:293:TRP:O	1:F:294:ASN:HB2	1.95	0.63
1:B:432:ILE:O	1:B:433:HIS:CG	2.51	0.63
1:E:311:PRO:O	1:E:313:ARG:HB2	1.98	0.63
1:B:479:LYS:O	1:B:480:SER:HB3	1.98	0.63
1:D:209:VAL:HG22	1:D:238:TYR:HB2	1.81	0.63
1:A:479:LYS:HD2	1:A:479:LYS:N	2.14	0.63
1:B:363:GLU:H	1:B:368:ALA:HA	1.64	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:460:ASP:O	1:D:462:LYS:N	2.30	0.62
1:D:311:PRO:CG	1:D:312:TRP:CD1	2.81	0.62
1:D:478:ASP:HA	1:D:479:LYS:HB3	1.81	0.62
1:F:364:ARG:C	1:F:366:GLY:HA3	2.19	0.62
1:B:469:GLY:HA2	1:B:470:GLU:HB2	1.80	0.62
1:F:458:HIS:HE1	1:F:471:GLU:HG2	1.65	0.62
1:B:67:ARG:HD2	1:B:291:ASP:HB2	1.82	0.62
1:E:293:TRP:O	1:E:351:ALA:HB3	2.00	0.62
1:B:216:LYS:HE3	1:B:216:LYS:O	1.99	0.62
1:D:431:ASN:HB3	1:D:433:HIS:N	2.13	0.62
1:A:310:GLU:HB3	1:A:311:PRO:CA	2.30	0.62
1:C:485:GLY:HA2	1:C:486:ILE:O	1.99	0.62
1:A:311:PRO:HA	1:A:313:ARG:HG3	1.81	0.62
1:D:469:GLY:CA	1:D:470:GLU:HB3	2.21	0.61
1:C:419:GLU:CB	1:C:420:LYS:HB3	2.29	0.61
1:B:487:LEU:HD13	1:B:501:ILE:HD11	1.83	0.61
1:E:367:GLY:CA	1:E:368:ALA:HB2	2.25	0.61
1:D:478:ASP:CB	1:D:479:LYS:HB3	2.31	0.61
1:C:145:PRO:HA	1:C:162:PRO:HG3	1.81	0.61
1:F:367:GLY:N	1:F:368:ALA:HB2	2.15	0.61
1:E:340:LYS:HG3	1:E:413:VAL:CG1	2.30	0.61
1:C:363:GLU:CB	1:C:368:ALA:HB3	2.27	0.61
1:F:46:ASP:HB3	1:F:52:LYS:HE3	1.82	0.61
1:B:362:THR:HA	1:B:369:ALA:H	1.66	0.60
1:D:477:SER:O	1:D:478:ASP:HB2	2.01	0.60
1:D:442:VAL:HG12	1:D:485:GLY:O	2.00	0.60
1:C:457:GLU:O	1:C:458:HIS:HB2	2.00	0.60
1:F:310:GLU:HB2	1:F:311:PRO:HA	1.84	0.60
1:F:364:ARG:O	1:F:366:GLY:HA3	2.02	0.60
1:F:298:HIS:HE1	1:F:327:ASP:OD2	1.85	0.60
1:B:491:LEU:HB3	1:B:492:ARG:CA	2.30	0.60
1:F:478:ASP:HA	1:F:479:LYS:CB	2.31	0.60
1:F:453:HIS:HD2	1:F:480:SER:HB2	1.66	0.60
1:C:420:LYS:H	1:C:421:GLU:CA	2.15	0.59
1:E:477:SER:O	1:E:478:ASP:HB2	2.01	0.59
1:E:312:TRP:HA	1:E:313:ARG:HB2	1.83	0.59
1:D:298:HIS:HE1	1:D:327:ASP:OD2	1.84	0.59
1:C:452:GLU:OE2	1:C:476:ASN:ND2	2.35	0.59
1:A:363:GLU:N	1:A:368:ALA:HB1	2.13	0.59
1:D:309:ASN:CA	1:D:310:GLU:HB2	2.33	0.58
1:B:317:PRO:HG3	1:B:364:ARG:HH21	1.67	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:17:ILE:HD12	1:F:389:VAL:HG23	1.85	0.58
1:B:188:TYR:HA	1:B:191:ILE:HG22	1.84	0.58
1:C:458:HIS:HA	1:C:459:GLN:CB	2.33	0.58
1:E:380:HIS:CD2	1:E:496:TRP:HE1	2.20	0.58
1:A:310:GLU:CB	1:A:311:PRO:CD	2.81	0.58
1:B:367:GLY:HA3	1:B:368:ALA:HB2	1.85	0.58
1:F:429:ASN:CA	1:F:430:ARG:O	2.43	0.58
1:F:7:THR:HG23	1:F:393:VAL:HG23	1.85	0.58
1:C:432:ILE:CG2	1:C:494:ALA:HB2	2.32	0.58
1:F:310:GLU:HB3	1:F:311:PRO:CA	2.34	0.57
1:D:363:GLU:O	1:D:364:ARG:HB2	2.04	0.57
1:C:385:GLY:HA2	1:C:415:ILE:HG12	1.86	0.57
1:A:458:HIS:HD2	1:A:460:ASP:H	1.52	0.57
1:E:380:HIS:HD2	1:E:496:TRP:HE1	1.51	0.57
1:B:336:ILE:HG23	1:B:413:VAL:HG22	1.85	0.57
1:B:491:LEU:HB3	1:B:492:ARG:C	2.25	0.57
1:D:482:PHE:CD1	1:D:486:ILE:O	2.57	0.57
1:C:485:GLY:HA2	1:C:486:ILE:C	2.25	0.57
1:A:312:TRP:H	1:A:313:ARG:HB2	1.68	0.57
1:D:307:MET:N	1:D:309:ASN:HB2	2.19	0.57
1:B:431:ASN:HA	1:B:432:ILE:CG2	2.30	0.57
1:B:51:ARG:HH22	1:B:368:ALA:CB	2.17	0.57
1:B:225:GLU:N	1:B:225:GLU:OE1	2.38	0.57
1:B:429:ASN:ND2	1:B:491:LEU:CD1	2.68	0.57
1:A:431:ASN:CA	1:A:432:ILE:O	2.53	0.56
1:A:410:ILE:HD12	1:A:491:LEU:HD12	1.87	0.56
1:B:348:ALA:HB1	1:B:349:CYS:CA	2.35	0.56
1:D:114:CYS:HA	1:D:117:VAL:HG22	1.86	0.56
1:E:308:GLN:H	1:E:309:ASN:C	2.09	0.56
1:A:209:VAL:HG22	1:A:238:TYR:HB2	1.88	0.56
1:F:257:ALA:HA	1:F:430:ARG:HH21	1.71	0.56
1:F:362:THR:HB	1:F:368:ALA:HB1	1.88	0.56
1:D:478:ASP:CA	1:D:479:LYS:HB3	2.36	0.56
1:E:240:SER:HA	1:E:289:SER:O	2.05	0.56
1:D:453:HIS:HE1	1:D:497:ASN:HD22	1.53	0.55
1:A:309:ASN:CB	1:A:310:GLU:CA	2.80	0.55
1:E:482:PHE:HD1	1:E:487:LEU:HB2	1.69	0.55
1:E:365:ASN:N	1:E:366:GLY:HA3	2.20	0.55
1:F:310:GLU:CB	1:F:311:PRO:CA	2.83	0.55
1:A:182:HIS:HB3	1:F:194:GLU:OE1	2.06	0.55
1:A:307:MET:HA	1:A:308:GLN:C	2.27	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:14:ILE:HG13	1:F:391:GLN:HG2	1.89	0.55
1:E:432:ILE:O	1:E:433:HIS:HB2	2.07	0.55
1:F:15:ALA:HB2	1:F:343:ASP:HB3	1.88	0.55
1:B:15:ALA:HB2	1:B:343:ASP:HB3	1.88	0.55
1:B:448:TYR:HD2	1:B:501:ILE:HG22	1.71	0.55
1:B:437:VAL:HA	1:B:490:MET:H	1.72	0.55
1:F:470:GLU:HG3	4:F:2030:HOH:O	2.06	0.54
1:B:469:GLY:CA	1:B:470:GLU:CB	2.81	0.54
1:A:125:VAL:HG12	1:A:138:LEU:HD23	1.88	0.54
1:C:82:GLY:O	1:C:106:GLY:HA3	2.07	0.54
1:D:172:ASN:O	1:D:173:ALA:C	2.46	0.54
1:B:431:ASN:H	1:B:431:ASN:HD22	1.55	0.54
1:B:290:PHE:O	1:B:348:ALA:CB	2.55	0.54
1:F:175:ASP:OD2	1:F:218:MET:HG3	2.07	0.54
1:B:348:ALA:CA	1:B:349:CYS:CB	2.77	0.54
1:B:21:ILE:HA	1:B:348:ALA:H	1.73	0.54
1:E:429:ASN:O	1:E:494:ALA:HA	2.08	0.54
1:C:12:TYR:CE1	1:C:391:GLN:HG3	2.43	0.54
1:E:67:ARG:NH2	1:E:172:ASN:OD1	2.41	0.54
1:A:311:PRO:HA	1:A:313:ARG:CG	2.38	0.54
1:B:396:SER:HB2	1:B:410:ILE:HG22	1.90	0.53
1:F:168:TRP:HB2	1:F:208:LEU:HD23	1.91	0.53
1:C:27:GLU:HG2	1:C:29:LEU:HB2	1.90	0.53
1:D:67:ARG:HD2	1:D:291:ASP:HB2	1.90	0.53
1:F:440:SER:HB2	1:F:487:LEU:HB3	1.90	0.53
1:E:431:ASN:ND2	1:E:431:ASN:H	2.06	0.53
1:E:431:ASN:HD22	1:E:431:ASN:N	2.01	0.53
1:B:307:MET:CG	1:B:308:GLN:N	2.61	0.53
1:A:367:GLY:CA	1:A:368:ALA:CB	2.62	0.53
1:B:12:TYR:CE2	1:C:12:TYR:CE2	2.97	0.53
1:E:393:VAL:HG13	1:E:394:ILE:H	1.73	0.53
1:B:438:LEU:N	1:B:489:SER:O	2.42	0.53
1:A:311:PRO:C	1:A:312:TRP:CD1	2.82	0.53
1:A:312:TRP:N	1:A:313:ARG:CB	2.71	0.52
1:A:168:TRP:HB2	1:A:208:LEU:HD23	1.90	0.52
1:A:241:LEU:C	1:A:242:HIS:HD2	2.13	0.52
1:A:310:GLU:CB	1:A:311:PRO:CB	2.66	0.52
1:D:350:LEU:CA	1:D:351:ALA:CB	2.85	0.52
1:F:61:LEU:HD22	1:F:378:PHE:CD1	2.45	0.52
1:C:460:ASP:CA	1:C:461:LEU:CB	2.85	0.52
1:B:456:LEU:HB3	1:B:496:TRP:HB3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:380:HIS:CD2	1:B:496:TRP:HE1	2.27	0.52
1:F:336:ILE:HG23	1:F:413:VAL:HG22	1.91	0.52
1:B:216:LYS:O	1:B:216:LYS:HE2	2.10	0.52
1:E:310:GLU:H	1:E:311:PRO:CA	2.16	0.52
1:D:478:ASP:CA	1:D:479:LYS:CB	2.88	0.52
1:C:410:ILE:HD11	1:C:436:ILE:CD1	2.40	0.52
1:B:469:GLY:CA	1:B:470:GLU:HB3	2.37	0.52
1:F:209:VAL:HG22	1:F:238:TYR:HB2	1.92	0.52
1:B:174:MET:HG3	1:B:179:GLN:HE22	1.75	0.52
1:D:336:ILE:HG23	1:D:413:VAL:CG2	2.39	0.52
1:F:363:GLU:HB2	1:F:368:ALA:HA	1.92	0.51
1:C:458:HIS:HD2	1:C:459:GLN:O	1.93	0.51
1:D:482:PHE:HD1	1:D:486:ILE:O	1.93	0.51
1:B:51:ARG:HH22	1:B:368:ALA:HB3	1.75	0.51
1:A:293:TRP:O	1:A:294:ASN:HB2	2.10	0.51
1:E:443:ARG:NH1	1:E:443:ARG:HB3	2.25	0.51
1:E:188:TYR:HA	1:E:191:ILE:HG22	1.91	0.51
1:E:478:ASP:HA	1:E:479:LYS:CB	2.14	0.51
1:F:362:THR:HA	1:F:369:ALA:H	1.75	0.51
1:D:311:PRO:HG2	1:D:312:TRP:HE1	1.72	0.51
1:D:480:SER:CB	1:D:481:SER:CA	2.79	0.51
1:A:12:TYR:OH	1:D:391:GLN:HG2	2.10	0.51
1:B:19:LYS:O	1:B:64:PRO:HG2	2.11	0.51
1:E:312:TRP:HA	1:E:313:ARG:CB	2.40	0.51
1:F:453:HIS:CD2	1:F:480:SER:HB2	2.45	0.51
1:C:460:ASP:CA	1:C:461:LEU:HB2	2.39	0.51
1:F:322:ILE:HG22	1:F:464:ARG:HD2	1.93	0.51
1:A:311:PRO:CA	1:A:313:ARG:HG3	2.41	0.51
1:D:471:GLU:O	1:D:472:VAL:CG2	2.53	0.51
1:E:209:VAL:HG22	1:E:238:TYR:HB2	1.91	0.51
1:D:430:ARG:HA	1:D:431:ASN:C	2.16	0.50
1:A:173:ALA:O	1:A:179:GLN:NE2	2.44	0.50
1:B:429:ASN:O	1:B:494:ALA:HA	2.12	0.50
1:F:13:LYS:HG2	1:F:388:ILE:HG21	1.93	0.50
1:C:460:ASP:O	1:C:463:ILE:HB	2.11	0.50
1:E:7:THR:H	1:E:393:VAL:HG12	1.77	0.50
1:E:389:VAL:HA	1:E:415:ILE:HG22	1.92	0.50
1:F:172:ASN:O	1:F:173:ALA:C	2.50	0.50
1:A:313:ARG:HD3	1:A:316:PRO:HB3	1.93	0.50
1:E:51:ARG:HH22	1:E:368:ALA:HA	1.76	0.50
1:B:307:MET:C	1:B:309:ASN:H	2.15	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:51:ARG:HH12	1:C:367:GLY:H	1.60	0.50
1:A:465:ASN:ND2	1:A:471:GLU:H	2.09	0.50
1:B:243:GLN:O	1:B:293:TRP:HA	2.12	0.50
1:A:309:ASN:HB3	1:A:310:GLU:HA	1.91	0.50
1:D:293:TRP:O	1:D:351:ALA:HB3	2.11	0.50
1:E:203:ASP:O	1:E:206:ILE:HG12	2.12	0.50
1:F:240:SER:HA	1:F:289:SER:O	2.11	0.50
1:C:5:ARG:HB2	1:C:395:ASN:HB3	1.94	0.50
1:A:240:SER:HA	1:A:289:SER:O	2.11	0.50
1:D:262:LEU:HD11	1:D:290:PHE:CZ	2.47	0.50
1:C:114:CYS:HA	1:C:117:VAL:HG22	1.94	0.50
1:C:380:HIS:HD2	1:C:496:TRP:HE1	1.60	0.50
1:A:409:ASP:O	1:A:430:ARG:O	2.29	0.49
1:F:399:HIS:HD2	1:F:409:ASP:OD1	1.95	0.49
1:A:463:ILE:HD13	1:A:464:ARG:H	1.77	0.49
1:F:17:ILE:HD12	1:F:389:VAL:CG2	2.42	0.49
1:D:409:ASP:O	1:D:430:ARG:O	2.29	0.49
1:B:444:GLY:HA2	1:B:445:MET:CB	2.41	0.49
1:E:99:SER:HA	1:E:312:TRP:HB3	1.94	0.49
1:A:306:ILE:HG22	1:A:307:MET:HG3	1.94	0.49
1:C:293:TRP:O	1:C:294:ASN:HB2	2.13	0.49
1:B:71:GLY:O	1:B:74:VAL:HG12	2.12	0.49
1:D:188:TYR:HA	1:D:191:ILE:HG22	1.95	0.49
1:D:311:PRO:HA	1:D:313:ARG:HG3	1.94	0.49
1:D:363:GLU:O	1:D:364:ARG:CB	2.60	0.49
1:D:24:SER:OG	1:D:353:LEU:HB2	2.13	0.49
1:C:47:GLU:H	1:C:47:GLU:CD	2.15	0.49
1:B:363:GLU:OE1	1:B:363:GLU:HA	2.13	0.49
1:A:267:ARG:HD3	4:A:2027:HOH:O	2.13	0.49
1:E:214:SER:O	1:E:242:HIS:HB2	2.13	0.49
1:E:29:LEU:HD13	1:E:356:VAL:HG11	1.95	0.49
1:E:350:LEU:CA	1:E:351:ALA:HB3	2.34	0.48
1:B:488:THR:O	1:B:488:THR:N	2.46	0.48
1:C:168:TRP:HB2	1:C:208:LEU:HD23	1.95	0.48
1:D:350:LEU:CA	1:D:351:ALA:HB3	2.29	0.48
1:B:482:PHE:HD1	1:B:487:LEU:HB2	1.78	0.48
1:B:310:GLU:CB	1:B:311:PRO:CA	2.91	0.48
1:A:291:ASP:O	1:A:292:GLU:HG3	2.13	0.48
1:A:232:ALA:O	1:A:236:VAL:HG22	2.13	0.48
1:B:432:ILE:O	1:B:433:HIS:CB	2.60	0.48
1:D:175:ASP:OD2	1:D:213:SER:HA	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:409:ASP:HA	1:E:430:ARG:O	2.12	0.48
1:E:410:ILE:HD11	1:E:491:LEU:HD12	1.96	0.48
1:B:367:GLY:CA	1:B:368:ALA:CB	2.86	0.48
1:D:472:VAL:O	1:D:473:TYR:CB	2.57	0.48
1:E:323:TYR:H	1:E:372:GLN:HE22	1.60	0.48
1:F:114:CYS:HA	1:F:117:VAL:HG22	1.95	0.48
1:B:106:GLY:O	1:B:108:ASN:N	2.47	0.48
1:D:471:GLU:HA	1:D:473:TYR:CD1	2.49	0.48
1:D:15:ALA:HB2	1:D:343:ASP:CB	2.44	0.48
1:B:444:GLY:CA	1:B:445:MET:HB2	2.44	0.48
1:B:166:LYS:HE3	4:B:2019:HOH:O	2.13	0.48
1:D:330:LEU:HA	1:D:333:LEU:HD12	1.95	0.48
1:C:29:LEU:HD13	1:C:356:VAL:HG11	1.96	0.48
1:D:6:MET:HB2	1:D:394:ILE:HD12	1.95	0.48
1:A:309:ASN:HB2	1:A:310:GLU:CB	2.44	0.47
1:D:173:ALA:O	1:D:179:GLN:NE2	2.47	0.47
1:F:410:ILE:HD12	1:F:491:LEU:CD1	2.44	0.47
1:B:479:LYS:HG3	1:B:490:MET:HG3	1.96	0.47
1:B:114:CYS:HA	1:B:117:VAL:HG22	1.96	0.47
1:F:380:HIS:HD2	1:F:496:TRP:NE1	2.08	0.47
1:D:399:HIS:HD2	1:D:409:ASP:OD1	1.98	0.47
1:A:492:ARG:HG3	1:A:493:ARG:N	2.29	0.47
1:F:458:HIS:CE1	1:F:471:GLU:HG2	2.47	0.47
1:D:154:ARG:HG2	1:D:159:VAL:HB	1.96	0.47
1:D:311:PRO:HD2	1:D:312:TRP:HD1	1.79	0.47
1:F:477:SER:O	1:F:478:ASP:HB2	2.15	0.47
1:B:172:ASN:O	1:B:173:ALA:C	2.52	0.47
1:B:432:ILE:HA	1:B:432:ILE:HD13	1.68	0.47
1:F:363:GLU:H	1:F:368:ALA:CB	2.28	0.47
1:C:380:HIS:CD2	1:C:496:TRP:HE1	2.33	0.47
1:B:244:TYR:OH	2:B:998:HRH:H5'1	2.15	0.47
1:E:45:SER:HA	1:E:50:PHE:O	2.14	0.47
1:C:238:TYR:HA	1:C:287:TYR:O	2.15	0.47
1:B:290:PHE:O	1:B:348:ALA:HB1	2.15	0.47
1:B:229:LEU:O	1:B:233:TYR:HB2	2.15	0.47
1:C:291:ASP:CA	1:C:349:CYS:HB3	2.41	0.47
1:C:229:LEU:O	1:C:233:TYR:HB2	2.14	0.47
1:A:336:ILE:HG23	1:A:413:VAL:HG22	1.96	0.47
1:C:291:ASP:O	1:C:292:GLU:HG3	2.14	0.47
1:B:209:VAL:HG22	1:B:238:TYR:HB2	1.96	0.47
1:B:480:SER:CA	1:B:481:SER:CB	2.89	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:478:ASP:HB3	1:F:479:LYS:C	2.36	0.46
1:F:172:ASN:ND2	1:F:179:GLN:OE1	2.46	0.46
1:F:455:VAL:HG23	1:F:497:ASN:ND2	2.31	0.46
1:A:29:LEU:HD13	1:A:356:VAL:HG11	1.95	0.46
1:D:473:TYR:CD2	1:D:473:TYR:O	2.68	0.46
1:A:460:ASP:O	1:A:461:LEU:CB	2.63	0.46
1:E:443:ARG:HA	1:E:444:GLY:O	2.14	0.46
1:F:142:CYS:HB3	1:F:165:ILE:HD12	1.97	0.46
1:C:241:LEU:HD22	1:C:290:PHE:CE2	2.51	0.46
1:B:61:LEU:HD22	1:B:378:PHE:HD2	1.79	0.46
1:B:431:ASN:HB2	1:B:432:ILE:CB	2.42	0.46
1:C:172:ASN:O	1:C:173:ALA:C	2.53	0.46
1:A:351:ALA:HA	1:A:352:GLN:HA	1.70	0.46
1:E:366:GLY:CA	1:E:367:GLY:O	2.62	0.46
1:D:311:PRO:HD2	1:D:312:TRP:CD1	2.51	0.46
1:C:72:ASN:ND2	1:C:179:GLN:OE1	2.48	0.46
1:D:453:HIS:CE1	1:D:497:ASN:HD22	2.32	0.46
1:E:358:ALA:O	1:E:372:GLN:HG3	2.16	0.46
1:F:365:ASN:N	1:F:366:GLY:HA3	2.28	0.46
1:E:432:ILE:O	1:E:433:HIS:CB	2.63	0.46
1:B:380:HIS:HD2	1:B:496:TRP:HE1	1.64	0.46
1:C:340:LYS:HD2	1:C:392:PRO:HG3	1.97	0.46
1:E:393:VAL:HG13	1:E:394:ILE:N	2.31	0.46
1:B:351:ALA:HA	1:B:352:GLN:HA	1.67	0.46
1:B:431:ASN:HB2	1:B:432:ILE:CA	2.46	0.46
1:A:492:ARG:HG3	1:A:493:ARG:H	1.81	0.46
1:A:428:VAL:HG12	1:A:429:ASN:N	2.31	0.46
1:D:167:VAL:HG22	1:D:207:GLU:HB2	1.98	0.46
1:B:168:TRP:HB2	1:B:208:LEU:HD23	1.97	0.46
1:E:368:ALA:HA	1:E:369:ALA:CB	2.31	0.46
1:B:290:PHE:O	1:B:348:ALA:HB2	2.16	0.46
1:E:348:ALA:HA	1:E:349:CYS:CB	2.23	0.45
1:D:309:ASN:N	1:D:310:GLU:CB	2.76	0.45
1:C:15:ALA:HB2	1:C:343:ASP:CB	2.44	0.45
1:D:478:ASP:CG	1:D:479:LYS:HB3	2.37	0.45
1:C:214:SER:O	1:C:242:HIS:HB2	2.16	0.45
1:F:29:LEU:HD21	1:F:319:LEU:HD13	1.97	0.45
1:A:308:GLN:O	1:A:309:ASN:C	2.53	0.45
1:A:310:GLU:CB	1:A:311:PRO:CA	2.93	0.45
1:C:432:ILE:HA	1:C:433:HIS:HA	1.69	0.45
1:F:478:ASP:HA	1:F:479:LYS:HB2	1.96	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:354:ILE:O	1:C:356:VAL:N	2.46	0.45
1:E:478:ASP:CA	1:E:479:LYS:CB	2.83	0.45
1:E:323:TYR:H	1:E:372:GLN:NE2	2.14	0.45
1:D:313:ARG:HB3	1:D:316:PRO:HG3	1.99	0.45
1:F:365:ASN:N	1:F:366:GLY:CA	2.80	0.45
1:C:262:LEU:HD22	1:C:334:MET:HG2	1.98	0.45
1:E:298:HIS:HE1	1:E:327:ASP:OD2	2.00	0.45
1:C:367:GLY:CA	1:C:368:ALA:HB3	2.46	0.45
1:F:125:VAL:HG12	1:F:138:LEU:HD23	1.99	0.45
1:A:91:ARG:HH22	1:A:314:ILE:HD11	1.81	0.45
1:F:356:VAL:O	1:F:357:ILE:C	2.55	0.45
1:B:431:ASN:N	1:B:431:ASN:ND2	2.63	0.45
1:A:12:TYR:CE1	1:A:391:GLN:HG3	2.51	0.45
1:F:203:ASP:O	1:F:206:ILE:HG12	2.17	0.45
1:F:307:MET:HA	1:F:308:GLN:HA	1.69	0.45
1:E:371:ARG:HD3	1:E:375:PHE:CD1	2.52	0.45
1:D:214:SER:O	1:D:242:HIS:HB2	2.17	0.45
1:D:229:LEU:O	1:D:233:TYR:HB2	2.17	0.45
1:A:47:GLU:CD	1:A:47:GLU:H	2.20	0.45
1:D:61:LEU:HD22	1:D:378:PHE:HD1	1.82	0.45
1:C:419:GLU:HB3	1:C:420:LYS:CB	2.38	0.45
1:F:61:LEU:HD22	1:F:378:PHE:HD1	1.82	0.45
1:C:203:ASP:O	1:C:206:ILE:HG12	2.17	0.44
1:E:9:ASP:HB3	1:E:12:TYR:HB3	1.97	0.44
1:E:106:GLY:O	1:E:108:ASN:N	2.50	0.44
1:D:38:TYR:CZ	1:D:40:PRO:HG3	2.52	0.44
1:B:111:ALA:HA	1:B:121:ILE:HD11	1.99	0.44
1:C:430:ARG:O	1:C:431:ASN:HB3	2.18	0.44
1:C:364:ARG:HA	1:C:365:ASN:HA	1.66	0.44
1:B:412:SER:HA	1:B:426:PHE:O	2.18	0.44
1:F:351:ALA:HA	1:F:352:GLN:HA	1.60	0.44
1:D:100:ILE:HG12	1:D:312:TRP:O	2.18	0.44
1:F:478:ASP:CB	1:F:479:LYS:HB2	2.47	0.44
1:E:351:ALA:N	1:E:352:GLN:CA	2.77	0.44
1:A:478:ASP:CG	1:A:479:LYS:H	2.21	0.44
1:A:188:TYR:HA	1:A:191:ILE:HG22	1.99	0.44
1:D:469:GLY:CA	1:D:470:GLU:CB	2.87	0.44
1:C:459:GLN:O	1:C:460:ASP:HB3	2.18	0.44
1:B:125:VAL:O	1:B:171:GLY:HA2	2.18	0.44
1:B:429:ASN:HD21	1:B:436:ILE:HD11	1.83	0.44
1:B:458:HIS:CD2	1:B:460:ASP:H	2.24	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:478:ASP:HB3	1:D:479:LYS:C	2.37	0.44
1:F:452:GLU:HG2	1:F:500:ARG:HH11	1.82	0.44
1:A:432:ILE:HA	1:A:433:HIS:HA	1.70	0.44
1:E:23:GLY:N	1:E:348:ALA:O	2.48	0.44
1:C:419:GLU:C	1:C:421:GLU:HA	2.33	0.44
1:B:440:SER:HB2	1:B:487:LEU:HB3	2.00	0.44
1:D:371:ARG:HD3	1:D:375:PHE:CE1	2.53	0.44
1:D:471:GLU:HA	1:D:473:TYR:HD1	1.82	0.43
1:A:431:ASN:CA	1:A:432:ILE:HG13	2.46	0.43
1:B:304:ALA:CA	1:B:305:ASN:CB	2.95	0.43
1:E:175:ASP:OD1	1:E:214:SER:OG	2.29	0.43
1:F:34:TYR:HB2	1:F:315:ALA:HB2	2.00	0.43
1:D:431:ASN:CB	1:D:432:ILE:CA	2.89	0.43
1:B:429:ASN:ND2	1:B:491:LEU:HD11	2.33	0.43
1:F:14:ILE:HG22	1:F:389:VAL:HB	2.00	0.43
1:E:153:MET:HG2	1:E:157:HIS:CE1	2.53	0.43
1:F:295:VAL:HG22	1:F:330:LEU:HD23	2.00	0.43
1:F:306:ILE:HG12	1:F:306:ILE:H	1.46	0.43
1:E:312:TRP:CA	1:E:313:ARG:HB2	2.48	0.43
1:B:154:ARG:HG2	1:B:159:VAL:HB	2.00	0.43
1:D:410:ILE:HD12	1:D:491:LEU:HD12	2.01	0.43
1:F:363:GLU:H	1:F:368:ALA:HB1	1.83	0.43
1:C:333:LEU:HD22	1:C:430:ARG:NH1	2.31	0.43
1:D:121:ILE:HD11	1:D:123:MET:SD	2.57	0.43
1:D:246:GLY:HA2	1:D:297:TYR:CD1	2.53	0.43
1:B:216:LYS:HB3	1:B:265:PHE:CE1	2.53	0.43
1:B:257:ALA:HA	1:B:430:ARG:HH21	1.83	0.43
1:B:63:VAL:O	1:B:119:ALA:HB2	2.19	0.43
1:A:242:HIS:N	1:A:242:HIS:CD2	2.87	0.43
1:C:84:GLY:HA2	1:C:157:HIS:CD2	2.54	0.43
1:A:51:ARG:O	1:A:55:ILE:HD12	2.19	0.43
1:B:159:VAL:HG11	1:B:163:HIS:CE1	2.54	0.43
1:D:492:ARG:O	1:D:493:ARG:C	2.57	0.43
1:D:102:PRO:HG3	1:D:314:ILE:HD12	2.00	0.43
1:D:432:ILE:HB	1:D:433:HIS:CD2	2.53	0.43
1:A:410:ILE:CD1	1:A:491:LEU:HD12	2.48	0.43
1:C:433:HIS:N	1:C:433:HIS:ND1	2.67	0.43
1:E:443:ARG:HH11	1:E:443:ARG:HB3	1.81	0.43
1:C:149:LYS:HD3	1:C:150:TYR:CE2	2.54	0.43
1:D:125:VAL:HG12	1:D:138:LEU:HD23	2.00	0.43
1:C:444:GLY:HA3	1:C:445:MET:HA	1.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:445:MET:CA	1:E:446:LYS:CB	2.88	0.43
1:D:336:ILE:HG23	1:D:413:VAL:HG22	2.01	0.43
1:C:298:HIS:HE1	1:C:327:ASP:OD2	2.02	0.43
1:F:45:SER:HA	1:F:50:PHE:O	2.19	0.42
1:C:308:GLN:O	1:C:309:ASN:CB	2.67	0.42
1:D:106:GLY:O	1:D:108:ASN:N	2.52	0.42
1:A:431:ASN:O	1:A:493:ARG:O	2.36	0.42
1:D:100:ILE:CG1	1:D:312:TRP:O	2.68	0.42
1:E:359:PRO:HA	1:E:372:GLN:HB2	2.01	0.42
1:B:429:ASN:HB2	1:B:491:LEU:HD13	2.01	0.42
1:F:179:GLN:HB3	1:F:179:GLN:HE21	1.49	0.42
1:C:407:VAL:HG13	1:C:430:ARG:HH21	1.85	0.42
1:B:446:LYS:HA	1:B:447:ASP:HA	1.82	0.42
1:E:376:TYR:CG	1:E:474:PRO:HD3	2.55	0.42
1:A:433:HIS:H	1:A:434:GLU:HB2	1.73	0.42
1:B:23:GLY:HA3	1:B:348:ALA:O	2.20	0.42
1:D:484:ASP:HA	1:D:485:GLY:HA2	1.45	0.42
1:D:433:HIS:ND1	1:D:433:HIS:N	2.68	0.42
1:A:310:GLU:CB	1:A:311:PRO:HD3	2.36	0.42
1:E:312:TRP:N	1:E:312:TRP:CD1	2.88	0.42
1:E:457:GLU:CB	1:E:458:HIS:HA	2.47	0.42
1:D:291:ASP:O	1:D:292:GLU:HG3	2.19	0.42
1:C:448:TYR:HD1	1:C:501:ILE:HG22	1.85	0.42
1:F:291:ASP:HA	1:F:349:CYS:HB2	2.02	0.42
1:A:312:TRP:N	1:A:312:TRP:CD1	2.88	0.42
1:A:312:TRP:CA	1:A:313:ARG:CB	2.96	0.42
1:B:311:PRO:HB2	1:B:312:TRP:H	1.63	0.42
1:F:35:ASP:OD1	1:F:364:ARG:O	2.36	0.42
1:E:456:LEU:HB2	1:E:496:TRP:HD1	1.85	0.42
1:A:241:LEU:C	1:A:242:HIS:CD2	2.93	0.42
1:E:241:LEU:C	1:E:242:HIS:HD2	2.23	0.42
1:B:5:ARG:HB2	1:B:395:ASN:HB3	2.02	0.42
1:D:17:ILE:HG13	1:D:389:VAL:HG23	2.01	0.42
1:B:142:CYS:HB3	1:B:165:ILE:HD12	2.01	0.42
1:C:440:SER:HB2	1:C:487:LEU:HB3	2.01	0.42
1:A:460:ASP:CB	1:A:463:ILE:HB	2.47	0.41
1:E:114:CYS:HA	1:E:117:VAL:HG22	2.01	0.41
1:C:240:SER:HA	1:C:289:SER:O	2.20	0.41
1:B:252:THR:HG21	1:B:461:LEU:HD13	2.02	0.41
1:C:386:ARG:O	1:C:417:ASN:HB2	2.20	0.41
1:B:7:THR:HG23	1:B:393:VAL:HB	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:432:ILE:HA	1:D:433:HIS:HA	1.62	0.41
1:C:367:GLY:HA3	1:C:368:ALA:O	2.19	0.41
1:B:291:ASP:HA	1:B:349:CYS:HB2	1.98	0.41
1:E:482:PHE:CD1	1:E:487:LEU:HB2	2.54	0.41
1:C:19:LYS:O	1:C:64:PRO:HG2	2.21	0.41
1:D:293:TRP:O	1:D:294:ASN:HB2	2.20	0.41
1:A:477:SER:CB	1:A:478:ASP:CA	2.87	0.41
1:E:336:ILE:HG23	1:E:413:VAL:HG22	2.01	0.41
1:C:172:ASN:HD22	1:C:179:GLN:NE2	2.12	0.41
1:F:17:ILE:CD1	1:F:389:VAL:CG2	2.98	0.41
1:C:206:ILE:HD12	1:C:208:LEU:HD21	2.01	0.41
1:E:4:ALA:HB2	1:E:397:PRO:HD2	2.02	0.41
1:F:10:LYS:HG2	1:F:10:LYS:H	1.69	0.41
1:C:419:GLU:O	1:C:421:GLU:HG3	2.21	0.41
1:A:428:VAL:CG1	1:A:429:ASN:N	2.83	0.41
1:D:309:ASN:CB	1:D:310:GLU:HB2	2.51	0.41
1:C:241:LEU:C	1:C:242:HIS:HD2	2.24	0.41
1:A:34:TYR:CZ	1:A:102:PRO:HG2	2.55	0.41
1:A:311:PRO:O	1:A:312:TRP:HD1	2.03	0.41
1:E:67:ARG:HD2	1:E:291:ASP:HB2	2.02	0.41
1:A:366:GLY:HA2	1:A:367:GLY:HA2	1.75	0.41
1:F:478:ASP:CA	1:F:479:LYS:CB	2.96	0.41
1:D:168:TRP:HB2	1:D:208:LEU:HD23	2.02	0.41
1:A:308:GLN:HB3	1:A:309:ASN:H	1.63	0.41
1:A:310:GLU:CG	1:A:311:PRO:HB3	2.49	0.41
1:D:238:TYR:HA	1:D:287:TYR:O	2.21	0.41
1:C:447:ASP:OD1	1:C:447:ASP:N	2.54	0.41
1:A:480:SER:HB3	1:A:489:SER:HA	2.03	0.41
1:E:308:GLN:CB	1:E:309:ASN:CG	2.89	0.41
1:F:455:VAL:HG23	1:F:497:ASN:HD22	1.85	0.41
1:C:448:TYR:CD1	1:C:501:ILE:HG22	2.56	0.41
1:A:174:MET:HE3	1:A:174:MET:HB2	1.96	0.41
1:C:23:GLY:HA2	1:C:63:VAL:HG13	2.03	0.41
1:C:70:GLY:HA2	1:C:71:GLY:HA3	1.88	0.41
1:A:410:ILE:HD11	1:A:436:ILE:CD1	2.32	0.41
1:E:166:LYS:HB3	4:E:2020:HOH:O	2.21	0.41
1:F:139:LEU:HD23	1:F:202:ILE:HB	2.01	0.41
1:C:483:ASP:HA	1:C:484:ASP:HA	1.87	0.41
1:B:431:ASN:HD22	1:B:431:ASN:N	2.19	0.40
1:B:304:ALA:N	1:B:305:ASN:HB2	2.36	0.40
1:D:416:TYR:HD1	1:D:423:VAL:HG22	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:125:VAL:O	1:E:171:GLY:HA2	2.21	0.40
1:C:27:GLU:HA	1:C:69:PRO:O	2.21	0.40
1:E:439:VAL:HG12	1:E:488:THR:HG22	2.02	0.40
1:C:126:ASN:C	1:C:126:ASN:OD1	2.59	0.40
1:B:479:LYS:HG3	1:B:490:MET:HB3	2.03	0.40
1:E:365:ASN:C	1:E:365:ASN:HD22	2.24	0.40
1:B:348:ALA:HA	1:B:349:CYS:CB	2.43	0.40
1:F:111:ALA:HA	1:F:121:ILE:HD11	2.04	0.40
1:B:51:ARG:HH22	1:B:368:ALA:HB1	1.85	0.40
1:A:252:THR:HG21	1:A:461:LEU:HD13	2.03	0.40
1:C:241:LEU:HD11	1:C:288:LEU:HD23	2.03	0.40
1:B:298:HIS:HE1	1:B:327:ASP:OD2	2.04	0.40
1:E:291:ASP:CA	1:E:349:CYS:HB3	2.38	0.40
1:A:409:ASP:HA	1:A:430:ARG:O	2.22	0.40
1:D:371:ARG:HD3	1:D:375:PHE:CD1	2.57	0.40
1:A:358:ALA:O	1:A:372:GLN:HG3	2.22	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	499/513 (97%)	445 (89%)	37 (7%)	17 (3%)	5	19
1	B	497/513 (97%)	435 (88%)	40 (8%)	22 (4%)	3	12
1	C	498/513 (97%)	446 (90%)	32 (6%)	20 (4%)	4	15
1	D	492/513 (96%)	428 (87%)	42 (8%)	22 (4%)	3	12
1	E	493/513 (96%)	436 (88%)	32 (6%)	25 (5%)	2	9
1	F	498/513 (97%)	445 (89%)	40 (8%)	13 (3%)	7	26
All	All	2977/3078 (97%)	2635 (88%)	223 (8%)	119 (4%)	4	15

All (119) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	62	ASN
1	A	309	ASN
1	A	310	GLU
1	A	311	PRO
1	A	364	ARG
1	A	368	ALA
1	A	432	ILE
1	A	434	GLU
1	A	470	GLU
1	A	493	ARG
1	B	217	ASP
1	B	304	ALA
1	B	307	MET
1	B	310	GLU
1	B	311	PRO
1	B	349	CYS
1	B	481	SER
1	C	309	ASN
1	C	310	GLU
1	C	349	CYS
1	C	458	HIS
1	C	461	LEU
1	C	478	ASP
1	C	481	SER
1	D	351	ALA
1	D	364	ARG
1	D	431	ASN
1	D	432	ILE
1	D	446	LYS
1	D	447	ASP
1	D	473	TYR
1	D	478	ASP
1	D	493	ARG
1	E	62	ASN
1	E	107	ILE
1	E	309	ASN
1	E	311	PRO
1	E	349	CYS
1	E	351	ALA
1	E	364	ARG
1	E	368	ALA
1	E	369	ALA

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Mol	Chain	Res	Type
1	E	433	HIS
1	E	434	GLU
1	F	310	GLU
1	F	430	ARG
1	F	470	GLU
1	A	445	MET
1	A	478	ASP
1	A	480	SER
1	A	481	SER
1	B	62	ASN
1	B	107	ILE
1	B	173	ALA
1	B	364	ARG
1	B	433	HIS
1	B	470	GLU
1	C	418	GLU
1	C	432	ILE
1	C	486	ILE
1	D	107	ILE
1	D	309	ASN
1	D	367	GLY
1	D	430	ARG
1	D	434	GLU
1	D	470	GLU
1	E	367	GLY
1	E	445	MET
1	E	478	ASP
1	E	479	LYS
1	F	294	ASN
1	F	368	ALA
1	A	430	ARG
1	A	484	ASP
1	C	297	TYR
1	C	355	ASN
1	C	493	ARG
1	D	173	ALA
1	D	472	VAL
1	D	479	LYS
1	E	471	GLU
1	F	357	ILE
1	F	478	ASP
1	B	305	ASN

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Mol	Chain	Res	Type
1	B	432	ILE
1	B	491	LEU
1	C	308	GLN
1	C	311	PRO
1	E	394	ILE
1	E	432	ILE
1	E	447	ASP
1	A	357	ILE
1	B	46	ASP
1	B	357	ILE
1	B	489	SER
1	C	431	ASN
1	C	460	ASP
1	D	357	ILE
1	D	461	LEU
1	E	297	TYR
1	E	310	GLU
1	E	355	ASN
1	F	355	ASN
1	F	449	ARG
1	B	480	SER
1	C	307	MET
1	C	357	ILE
1	C	419	GLU
1	D	310	GLU
1	E	469	GLY
1	F	62	ASN
1	F	173	ALA
1	F	307	MET
1	B	367	GLY
1	B	306	ILE
1	E	317	PRO
1	E	357	ILE
1	D	311	PRO
1	F	311	PRO

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	434/450 (96%)	406 (94%)	28 (6%)	21	52
1	B	434/450 (96%)	398 (92%)	36 (8%)	14	38
1	C	428/450 (95%)	395 (92%)	33 (8%)	16	42
1	D	420/450 (93%)	393 (94%)	27 (6%)	22	53
1	E	417/450 (93%)	391 (94%)	26 (6%)	23	55
1	F	423/450 (94%)	396 (94%)	27 (6%)	22	53
All	All	2556/2700 (95%)	2379 (93%)	177 (7%)	19	48

All (177) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	LYS
1	A	11	ASP
1	A	12	TYR
1	A	13	LYS
1	A	98	LYS
1	A	112	LYS
1	A	132	ILE
1	A	241	LEU
1	A	290	PHE
1	A	302	GLU
1	A	307	MET
1	A	312	TRP
1	A	365	ASN
1	A	390	LEU
1	A	413	VAL
1	A	415	ILE
1	A	418	GLU
1	A	430	ARG
1	A	432	ILE
1	A	436	ILE
1	A	445	MET
1	A	446	LYS
1	A	447	ASP
1	A	449	ARG
1	A	455	VAL
1	A	463	ILE
1	A	479	LYS
1	A	488	THR

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Mol	Chain	Res	Type
1	B	11	ASP
1	B	12	TYR
1	B	17	ILE
1	B	43	SER
1	B	87	GLU
1	B	98	LYS
1	B	174	MET
1	B	179	GLN
1	B	216	LYS
1	B	243	GLN
1	B	290	PHE
1	B	305	ASN
1	B	310	GLU
1	B	364	ARG
1	B	365	ASN
1	B	380	HIS
1	B	394	ILE
1	B	413	VAL
1	B	419	GLU
1	B	429	ASN
1	B	431	ASN
1	B	432	ILE
1	B	435	ASP
1	B	445	MET
1	B	446	LYS
1	B	447	ASP
1	B	455	VAL
1	B	463	ILE
1	B	468	ASN
1	B	476	ASN
1	B	480	SER
1	B	483	ASP
1	B	487	LEU
1	B	489	SER
1	B	490	MET
1	B	491	LEU
1	C	7	THR
1	C	12	TYR
1	C	13	LYS
1	C	16	GLU
1	C	17	ILE
1	C	43	SER

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Mol	Chain	Res	Type
1	C	44	LYS
1	C	47	GLU
1	C	56	GLU
1	C	98	LYS
1	C	112	LYS
1	C	134	ASP
1	C	160	LYS
1	C	218	MET
1	C	241	LEU
1	C	250	ASN
1	C	307	MET
1	C	365	ASN
1	C	390	LEU
1	C	415	ILE
1	C	419	GLU
1	C	429	ASN
1	C	430	ARG
1	C	432	ILE
1	C	433	HIS
1	C	436	ILE
1	C	458	HIS
1	C	460	ASP
1	C	463	ILE
1	C	468	ASN
1	C	479	LYS
1	C	480	SER
1	C	483	ASP
1	D	16	GLU
1	D	98	LYS
1	D	121	ILE
1	D	160	LYS
1	D	220	THR
1	D	290	PHE
1	D	308	GLN
1	D	314	ILE
1	D	350	LEU
1	D	393	VAL
1	D	394	ILE
1	D	413	VAL
1	D	415	ILE
1	D	418	GLU
1	D	430	ARG

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Mol	Chain	Res	Type
1	D	433	HIS
1	D	436	ILE
1	D	440	SER
1	D	445	MET
1	D	455	VAL
1	D	459	GLN
1	D	463	ILE
1	D	464	ARG
1	D	468	ASN
1	D	478	ASP
1	D	488	THR
1	D	490	MET
1	E	44	LYS
1	E	47	GLU
1	E	87	GLU
1	E	98	LYS
1	E	107	ILE
1	E	160	LYS
1	E	179	GLN
1	E	290	PHE
1	E	305	ASN
1	E	312	TRP
1	E	349	CYS
1	E	350	LEU
1	E	365	ASN
1	E	372	GLN
1	E	390	LEU
1	E	391	GLN
1	E	410	ILE
1	E	413	VAL
1	E	431	ASN
1	E	443	ARG
1	E	448	TYR
1	E	456	LEU
1	E	457	GLU
1	E	463	ILE
1	E	486	ILE
1	E	487	LEU
1	F	10	LYS
1	F	11	ASP
1	F	14	ILE
1	F	17	ILE

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Mol	Chain	Res	Type
1	F	44	LYS
1	F	47	GLU
1	F	98	LYS
1	F	107	ILE
1	F	160	LYS
1	F	179	GLN
1	F	242	HIS
1	F	283	LYS
1	F	290	PHE
1	F	306	ILE
1	F	365	ASN
1	F	388	ILE
1	F	390	LEU
1	F	393	VAL
1	F	413	VAL
1	F	415	ILE
1	F	421	GLU
1	F	430	ARG
1	F	446	LYS
1	F	449	ARG
1	F	453	HIS
1	F	455	VAL
1	F	480	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (52) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	242	HIS
1	A	243	GLN
1	A	298	HIS
1	A	429	ASN
1	A	458	HIS
1	A	465	ASN
1	B	72	ASN
1	B	104	GLN
1	B	243	GLN
1	B	298	HIS
1	B	380	HIS
1	B	395	ASN
1	B	399	HIS
1	B	429	ASN
1	B	458	HIS

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Mol	Chain	Res	Type
1	B	476	ASN
1	B	497	ASN
1	C	72	ASN
1	C	104	GLN
1	C	179	GLN
1	C	298	HIS
1	C	380	HIS
1	C	429	ASN
1	C	458	HIS
1	D	72	ASN
1	D	104	GLN
1	D	242	HIS
1	D	298	HIS
1	D	305	ASN
1	D	308	GLN
1	D	399	HIS
1	D	429	ASN
1	D	453	HIS
1	D	458	HIS
1	E	72	ASN
1	E	179	GLN
1	E	242	HIS
1	E	298	HIS
1	E	365	ASN
1	E	380	HIS
1	E	429	ASN
1	E	431	ASN
1	E	459	GLN
1	E	476	ASN
1	F	72	ASN
1	F	137	ASN
1	F	298	HIS
1	F	380	HIS
1	F	399	HIS
1	F	453	HIS
1	F	458	HIS
1	F	497	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 ⓘ

12 carbohydrates 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	AHR	A	998	2	9,9,10	0.49	0	12,12,14	1.20	1 (8%)
2	XYS	A	999	2	9,9,10	0.85	0	12,12,14	1.84	3 (25%)
2	AHR	B	998	2	9,9,10	0.54	0	12,12,14	1.57	3 (25%)
2	XYS	B	999	2	9,9,10	0.92	0	12,12,14	1.94	3 (25%)
2	AHR	C	998	2	9,9,10	0.41	0	12,12,14	0.92	1 (8%)
2	XYS	C	999	2	9,9,10	0.81	0	12,12,14	2.07	3 (25%)
2	AHR	D	998	2	9,9,10	0.51	0	12,12,14	1.09	1 (8%)
2	XYS	D	999	2	9,9,10	0.96	0	12,12,14	2.11	4 (33%)
2	AHR	E	998	2	9,9,10	0.57	0	12,12,14	1.13	1 (8%)
2	XYS	E	999	2	9,9,10	0.91	0	12,12,14	1.57	2 (16%)
2	AHR	F	998	2	9,9,10	0.49	0	12,12,14	1.00	1 (8%)
2	XYS	F	999	2	9,9,10	0.91	0	12,12,14	1.49	2 (16%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	AHR	A	998	2	-	0/2/15/18	0/1/1/1
2	XYS	A	999	2	-	0/0/14/17	1/1/1/1
2	AHR	B	998	2	-	0/2/15/18	0/1/1/1
2	XYS	B	999	2	-	0/0/14/17	1/1/1/1
2	AHR	C	998	2	-	0/2/15/18	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	XYS	C	999	2	-	0/0/14/17	0/1/1/1
2	AHR	D	998	2	-	0/2/15/18	0/1/1/1
2	XYS	D	999	2	-	0/0/14/17	1/1/1/1
2	AHR	E	998	2	-	0/2/15/18	0/1/1/1
2	XYS	E	999	2	-	0/0/14/17	0/1/1/1
2	AHR	F	998	2	-	0/2/15/18	0/1/1/1
2	XYS	F	999	2	-	0/0/14/17	0/1/1/1

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	999	XYS	C4-C3-C2	-3.34	107.94	111.24
2	D	999	XYS	C4-C3-C2	-3.34	107.94	111.24
2	B	999	XYS	C4-C3-C2	-2.85	108.42	111.24
2	A	999	XYS	C4-C3-C2	-2.70	108.57	111.24
2	B	998	AHR	C5'-C4'-C3'	-2.43	109.34	115.08
2	E	999	XYS	O4-C4-C5	-2.32	104.55	109.21
2	D	999	XYS	O3-C3-C2	2.01	113.63	110.00
2	F	999	XYS	C5-O5-C1	2.22	115.14	111.57
2	F	998	AHR	C1'-C2'-C3'	2.44	105.57	101.64
2	E	998	AHR	C1'-C2'-C3'	2.48	105.63	101.64
2	B	998	AHR	O4'-C4'-C3'	2.49	108.82	104.43
2	C	998	AHR	C1'-C2'-C3'	2.61	105.84	101.64
2	B	999	XYS	C5-O5-C1	2.74	115.98	111.57
2	A	998	AHR	C1'-C2'-C3'	2.93	106.36	101.64
2	D	998	AHR	C1'-C2'-C3'	2.97	106.43	101.64
2	C	999	XYS	C5-O5-C1	3.13	116.62	111.57
2	D	999	XYS	C5-O5-C1	3.26	116.83	111.57
2	A	999	XYS	C5-O5-C1	3.32	116.92	111.57
2	B	998	AHR	C1'-C2'-C3'	3.44	107.18	101.64
2	A	999	XYS	O3-C3-C4	3.74	116.75	110.00
2	E	999	XYS	O3-C3-C4	3.92	117.09	110.00
2	F	999	XYS	O3-C3-C4	4.02	117.26	110.00
2	D	999	XYS	O3-C3-C4	4.67	118.44	110.00
2	B	999	XYS	O3-C3-C4	4.87	118.80	110.00
2	C	999	XYS	O3-C3-C4	5.32	119.60	110.00

There are no chirality outliers.

There are no torsion outliers.

All (3) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	999	XYS	C1-C2-C3-C4-C5-O5
2	A	999	XYS	C1-C2-C3-C4-C5-O5
2	D	999	XYS	C1-C2-C3-C4-C5-O5

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	998	AHR	1	0

5.6 Ligand geometry

9 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	$\# Z > 2$
3	EDO	A	1503	-	3,3,3	0.56	0	2,2,2	0.22	0
3	EDO	A	1504	-	3,3,3	0.54	0	2,2,2	0.27	0
3	EDO	B	1503	-	3,3,3	0.58	0	2,2,2	0.19	0
3	EDO	C	1503	-	3,3,3	0.55	0	2,2,2	0.28	0
3	EDO	C	1504	-	3,3,3	0.49	0	2,2,2	0.42	0
3	EDO	D	1503	-	3,3,3	0.48	0	2,2,2	0.38	0
3	EDO	E	1503	-	3,3,3	0.65	0	2,2,2	0.27	0
3	EDO	E	1504	-	3,3,3	0.44	0	2,2,2	0.57	0
3	EDO	F	1503	-	3,3,3	0.48	0	2,2,2	0.32	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	A	1503	-	-	0/1/1/1	0/0/0/0
3	EDO	A	1504	-	-	0/1/1/1	0/0/0/0
3	EDO	B	1503	-	-	0/1/1/1	0/0/0/0
3	EDO	C	1503	-	-	0/1/1/1	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	C	1504	-	-	0/1/1/1	0/0/0/0
3	EDO	D	1503	-	-	0/1/1/1	0/0/0/0
3	EDO	E	1503	-	-	0/1/1/1	0/0/0/0
3	EDO	E	1504	-	-	0/1/1/1	0/0/0/0
3	EDO	F	1503	-	-	0/1/1/1	0/0/0/0

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.

No monomer is involved in short contacts.

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	501/513 (97%)	0.08	9 (1%) 71 68	49, 53, 56, 64	0
1	B	500/513 (97%)	-0.01	7 (1%) 78 76	49, 53, 57, 64	0
1	C	499/513 (97%)	0.09	10 (2%) 68 64	49, 53, 57, 64	0
1	D	498/513 (97%)	0.04	4 (0%) 87 86	49, 53, 56, 64	0
1	E	497/513 (96%)	0.13	12 (2%) 62 57	49, 53, 56, 62	0
1	F	500/513 (97%)	0.11	8 (1%) 74 72	49, 53, 57, 64	0
All	All	2995/3078 (97%)	0.07	50 (1%) 73 70	49, 53, 56, 64	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	501	ILE	4.1
1	A	501	ILE	3.9
1	D	437	VAL	3.6
1	D	431	ASN	3.2
1	C	484	ASP	3.1
1	A	448	TYR	3.1
1	F	448	TYR	2.9
1	B	445	MET	2.9
1	E	398	LEU	2.8
1	E	498	VAL	2.8
1	F	481	SER	2.8
1	E	490	MET	2.8
1	C	307	MET	2.7
1	E	439	VAL	2.7
1	B	446	LYS	2.6
1	A	476	ASN	2.6
1	E	431	ASN	2.6
1	F	447	ASP	2.5
1	A	485	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	398	LEU	2.5
1	D	433	HIS	2.4
1	E	478	ASP	2.4
1	C	420	LYS	2.4
1	C	437	VAL	2.4
1	C	486	ILE	2.4
1	E	436	ILE	2.3
1	F	430	ARG	2.3
1	A	441	ASP	2.3
1	B	441	ASP	2.3
1	E	440	SER	2.3
1	B	447	ASP	2.3
1	F	431	ASN	2.3
1	A	443	ARG	2.3
1	E	4	ALA	2.3
1	B	429	ASN	2.3
1	F	443	ARG	2.2
1	C	478	ASP	2.2
1	A	484	ASP	2.2
1	E	457	GLU	2.1
1	A	447	ASP	2.1
1	E	50	PHE	2.1
1	B	416	TYR	2.1
1	D	429	ASN	2.1
1	F	429	ASN	2.1
1	E	220	THR	2.1
1	C	446	LYS	2.1
1	C	416	TYR	2.1
1	B	437	VAL	2.0
1	C	490	MET	2.0
1	F	367	GLY	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](#)

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	XYS	F	999	9/10	0.94	0.21	0.79	58,60,61,61	0
2	XYS	E	999	9/10	0.92	0.21	0.68	58,60,61,61	0
2	AHR	D	998	9/10	0.95	0.19	0.24	54,55,57,57	0
2	AHR	A	998	9/10	0.92	0.21	0.11	54,56,57,57	0
2	AHR	B	998	9/10	0.96	0.19	0.07	54,55,57,57	0
2	XYS	A	999	9/10	0.88	0.19	-0.55	58,60,61,61	0
2	AHR	E	998	9/10	0.95	0.17	-0.61	54,55,57,57	0
2	XYS	B	999	9/10	0.93	0.17	-0.75	58,60,61,61	0
2	XYS	D	999	9/10	0.93	0.16	-1.08	58,60,61,61	0
2	AHR	C	998	9/10	0.94	0.18	-1.19	54,56,57,57	0
2	XYS	C	999	9/10	0.94	0.17	-1.61	58,60,61,61	0
2	AHR	F	998	9/10	0.94	0.13	-2.07	54,55,56,57	0

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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
3	EDO	A	1503	4/4	0.83	0.40	11.36	55,55,55,56	0
3	EDO	E	1503	4/4	0.89	0.35	4.59	55,56,56,57	0
3	EDO	C	1503	4/4	0.78	0.29	4.36	57,59,59,60	0
3	EDO	F	1503	4/4	0.87	0.28	3.67	56,56,56,57	0
3	EDO	B	1503	4/4	0.66	0.25	2.00	56,57,58,58	0
3	EDO	D	1503	4/4	0.80	0.18	0.11	57,57,57,58	0
3	EDO	C	1504	4/4	0.87	0.16	-0.73	52,53,55,56	0
3	EDO	E	1504	4/4	0.90	0.15	-0.97	57,58,59,60	0
3	EDO	A	1504	4/4	0.88	0.17	-2.40	53,54,54,54	0

6.5 Other polymers ⓘ

There are no such residues in this entry.