



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 07:09 PM GMT

PDB ID : 1AFC
Title : STRUCTURAL STUDIES OF THE BINDING OF THE ANTI-ULCER
DRUG SUCROSE OCTASULFATE TO ACIDIC FIBROBLAST GROWTH
FACTOR
Authors : Zhu, X.; Hsu, B.T.; Rees, D.C.
Deposited on : 1993-07-13
Resolution : 2.70 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

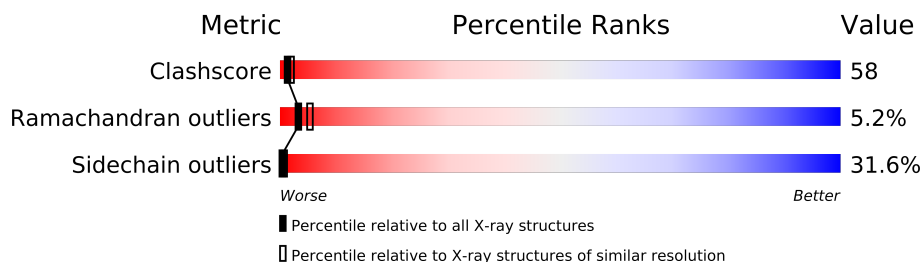
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : NOT EXECUTED
EDS : NOT EXECUTED
Percentile statistics : 21963
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	140	
1	B	140	
1	C	140	
1	D	140	
1	E	140	
1	F	140	
1	G	140	
1	H	140	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8304 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

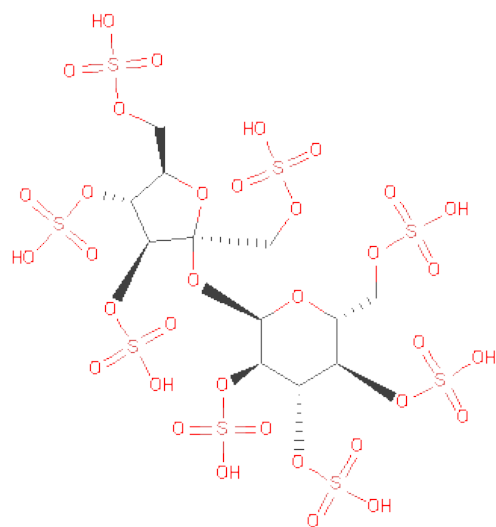
- Molecule 1 is a protein called ACIDIC FIBROBLAST GROWTH FACTOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	127	Total	C	N	O	S	0	0	0
			983	630	168	182	3			
1	B	127	Total	C	N	O	S	0	0	0
			983	630	168	182	3			
1	C	127	Total	C	N	O	S	0	0	0
			983	630	168	182	3			
1	D	127	Total	C	N	O	S	0	0	0
			983	630	168	182	3			
1	E	127	Total	C	N	O	S	0	0	0
			983	630	168	182	3			
1	F	127	Total	C	N	O	S	0	0	0
			983	630	168	182	3			
1	G	127	Total	C	N	O	S	0	0	0
			983	630	168	182	3			
1	H	127	Total	C	N	O	S	0	0	0
			983	630	168	182	3			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	47	ALA	CYS	CONFLICT	UNP P03968
B	47	ALA	CYS	CONFLICT	UNP P03968
C	47	ALA	CYS	CONFLICT	UNP P03968
D	47	ALA	CYS	CONFLICT	UNP P03968
E	47	ALA	CYS	CONFLICT	UNP P03968
F	47	ALA	CYS	CONFLICT	UNP P03968
G	47	ALA	CYS	CONFLICT	UNP P03968
H	47	ALA	CYS	CONFLICT	UNP P03968

- Molecule 2 is SUGAR (SUCROSE OCTASULFATE) (three-letter code: SCR) (formula: $C_{12}H_{22}O_{35}S_8$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	S	0	0
			55	12	35	8		
2	B	1	Total	C	O	S	0	0
			55	12	35	8		
2	C	1	Total	C	O	S	0	0
			55	12	35	8		
2	D	1	Total	C	O	S	0	0
			55	12	35	8		
2	E	1	Total	C	O	S	0	0
			55	12	35	8		
2	F	1	Total	C	O	S	0	0
			55	12	35	8		
2	G	1	Total	C	O	S	0	0
			55	12	35	8		
2	H	1	Total	C	O	S	0	0
			55	12	35	8		

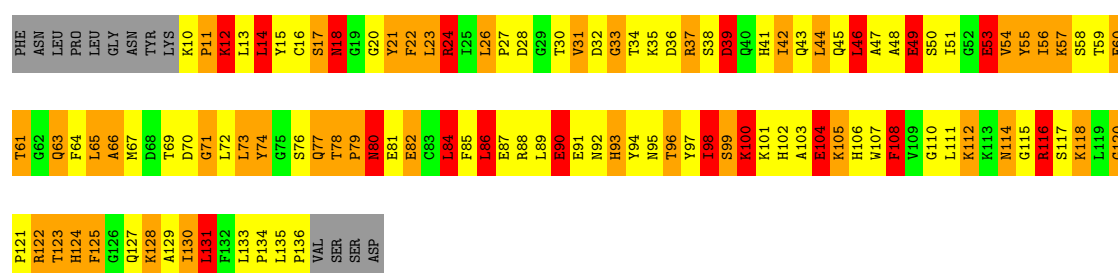
3 Residue-property plots

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.

Note EDS was not executed.

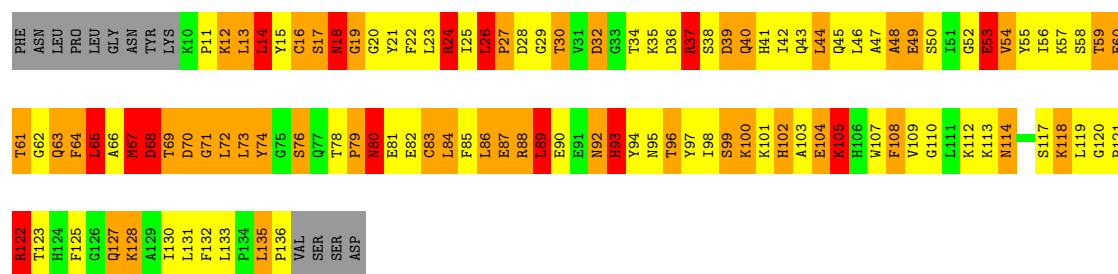
• Molecule 1: ACIDIC FIBROBLAST GROWTH FACTOR

Chain A:



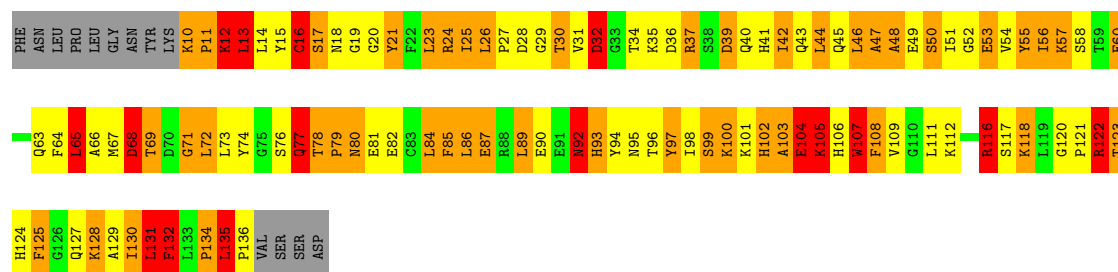
• Molecule 1: ACIDIC FIBROBLAST GROWTH FACTOR

Chain B:



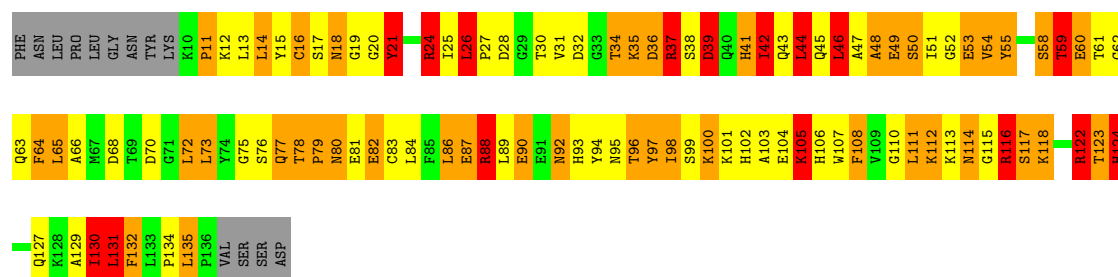
• Molecule 1: ACIDIC FIBROBLAST GROWTH FACTOR

Chain C:



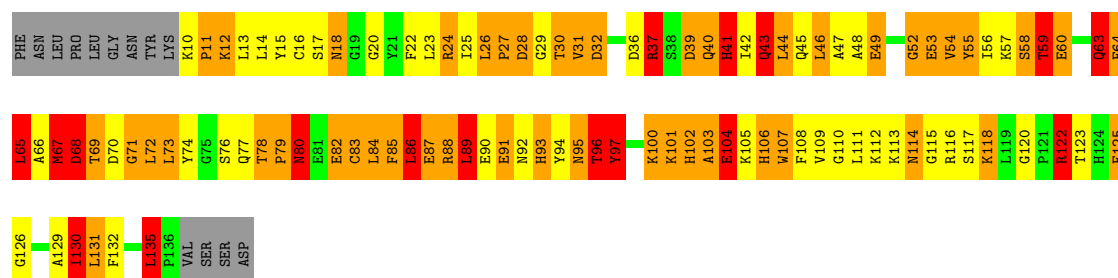
• Molecule 1: ACIDIC FIBROBLAST GROWTH FACTOR

Chain D:



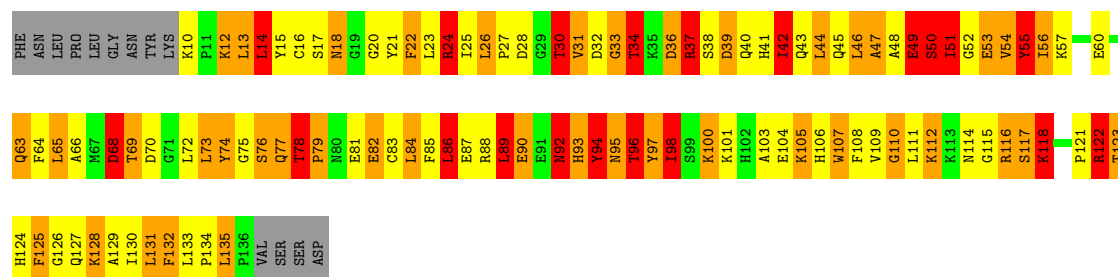
- Molecule 1: ACIDIC FIBROBLAST GROWTH FACTOR

Chain E:



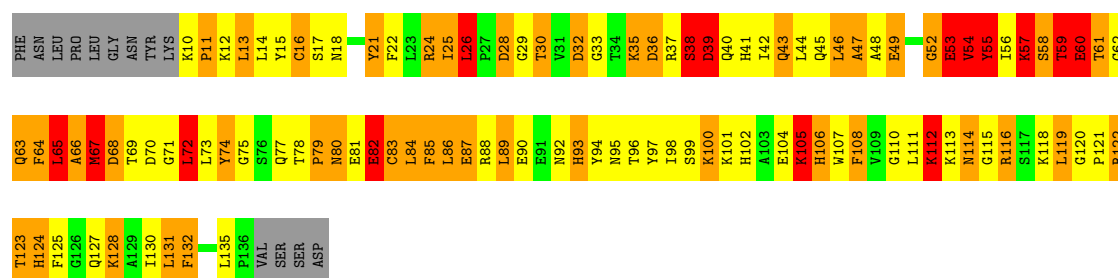
- Molecule 1: ACIDIC FIBROBLAST GROWTH FACTOR

Chain F:



- Molecule 1: ACIDIC FIBROBLAST GROWTH FACTOR

Chain G:



- Molecule 1: ACIDIC FIBROBLAST GROWTH FACTOR

Chain H:



R122	G62	PHE
T123	Q63	ASN
H124	F64	LEU
F125	G65	PRO
G126	A66	LEU
Q127	M67	GLY
K128	D68	ASN
A129	T69	TYR
I130	D70	LVS
L131	G71	K10
F132	L72	P11
L133	L73	K12
P134	Y74	L13
L135	G75	L14
P136	S76	
VAL	Q77	S17
SER	T78	N18
SER	P79	G19
ASP	R80	G20
	E81	Y21
	E82	F22
	C83	L23
	L84	R24
	F85	T25
	L86	L26
	E87	P27
	R88	D28
	L89	G29
	E90	T30
	E91	V31
	N92	D32
	H93	G33
	Y94	T34
	N95	K35
	T96	D36
	Y97	R37
	I98	S38
	S99	D39
	K100	Q40
	K101	H41
	H102	L42
	A103	Q43
	E104	L44
	K105	Q45
	H106	L46
	V107	A47
	F108	A48
	V109	E49
	G110	S50
	L111	I51
	K112	G52
	K113	E53
	N114	V54
	G115	Y55
	R116	I56
	S117	K57
	K118	S58
	L119	T59
	G120	E60
	P121	T61

4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	110.60Å 110.60Å 172.20Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	(Not available) – 2.70	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.70)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.204 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	8304	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.87	0/1007	1.54	8/1363 (0.6%)
1	B	0.91	0/1007	1.48	8/1363 (0.6%)
1	C	0.89	0/1007	1.50	10/1363 (0.7%)
1	D	0.89	0/1007	1.49	11/1363 (0.8%)
1	E	0.91	0/1007	1.49	9/1363 (0.7%)
1	F	0.90	0/1007	1.52	8/1363 (0.6%)
1	G	0.89	0/1007	1.46	4/1363 (0.3%)
1	H	0.90	0/1007	1.51	8/1363 (0.6%)
All	All	0.89	0/8056	1.50	66/10904 (0.6%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	89
1	B	0	77
1	C	0	76
1	D	0	76
1	E	0	74
1	F	0	78
1	G	0	67
1	H	0	83
All	All	0	620

There are no bond length outliers.

All (66) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	107	TRP	CD1-CG-CD2	11.01	115.11	106.30
1	A	24	ARG	NE-CZ-NH2	-10.65	114.97	120.30
1	D	37	ARG	NE-CZ-NH2	-10.06	115.27	120.30
1	E	24	ARG	NE-CZ-NH2	-9.73	115.43	120.30
1	D	24	ARG	NE-CZ-NH2	-9.52	115.54	120.30
1	A	107	TRP	CD1-CG-CD2	9.20	113.66	106.30
1	C	107	TRP	CD1-CG-CD2	8.88	113.40	106.30
1	F	24	ARG	NE-CZ-NH1	8.70	124.65	120.30
1	H	122	ARG	NE-CZ-NH1	-8.62	115.99	120.30
1	E	107	TRP	CD1-CG-CD2	8.08	112.76	106.30
1	F	107	TRP	CE2-CD2-CG	-8.02	100.88	107.30
1	D	107	TRP	CD1-CG-CD2	7.98	112.68	106.30
1	B	107	TRP	CD1-CG-CD2	7.89	112.62	106.30
1	F	122	ARG	NE-CZ-NH2	-7.63	116.48	120.30
1	E	24	ARG	NE-CZ-NH1	7.53	124.07	120.30
1	A	107	TRP	CE2-CD2-CG	-7.48	101.31	107.30
1	H	107	TRP	CD1-CG-CD2	7.47	112.27	106.30
1	B	122	ARG	NE-CZ-NH2	-7.46	116.57	120.30
1	C	107	TRP	CE2-CD2-CG	-7.46	101.33	107.30
1	A	21	TYR	O-C-N	-7.37	110.91	122.70
1	G	107	TRP	CD1-CG-CD2	7.29	112.13	106.30
1	G	55	TYR	CB-CG-CD2	-7.22	116.67	121.00
1	F	107	TRP	CG-CD1-NE1	-7.04	103.06	110.10
1	D	39	ASP	CB-CG-OD2	6.91	124.52	118.30
1	B	107	TRP	CE2-CD2-CG	-6.79	101.87	107.30
1	E	107	TRP	CE2-CD2-CG	-6.72	101.92	107.30
1	D	116	ARG	NE-CZ-NH2	-6.67	116.97	120.30
1	D	107	TRP	CE2-CD2-CG	-6.49	102.11	107.30
1	E	37	ARG	NE-CZ-NH2	-6.46	117.07	120.30
1	C	18	ASN	C-N-CA	-6.42	108.81	122.30
1	H	107	TRP	CE2-CD2-CG	-6.42	102.16	107.30
1	C	32	ASP	CB-CG-OD2	-6.24	112.68	118.30
1	A	125	PHE	CB-CG-CD1	-6.21	116.46	120.80
1	G	107	TRP	CE2-CD2-CG	-6.17	102.37	107.30
1	D	122	ARG	CD-NE-CZ	-6.15	114.99	123.60
1	C	24	ARG	NE-CZ-NH1	6.13	123.36	120.30
1	B	68	ASP	CB-CG-OD1	-5.99	112.91	118.30
1	B	24	ARG	NE-CZ-NH2	-5.98	117.31	120.30
1	A	107	TRP	CB-CG-CD1	-5.91	119.32	127.00
1	H	48	ALA	CB-CA-C	-5.85	101.32	110.10
1	E	78	THR	N-CA-CB	-5.75	99.38	110.30
1	D	21	TYR	CB-CG-CD2	-5.61	117.64	121.00
1	C	122	ARG	NE-CZ-NH1	-5.51	117.55	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	74	TYR	CB-CG-CD2	-5.50	117.70	121.00
1	A	107	TRP	CG-CD1-NE1	-5.50	104.60	110.10
1	D	116	ARG	NE-CZ-NH1	5.49	123.04	120.30
1	H	55	TYR	CB-CG-CD2	-5.46	117.72	121.00
1	H	27	PRO	O-C-N	-5.46	113.96	122.70
1	D	122	ARG	NE-CZ-NH2	-5.45	117.57	120.30
1	B	107	TRP	CG-CD1-NE1	-5.45	104.65	110.10
1	C	107	TRP	CG-CD1-NE1	-5.43	104.67	110.10
1	A	24	ARG	NE-CZ-NH1	5.42	123.01	120.30
1	F	55	TYR	CB-CG-CD2	-5.42	117.75	121.00
1	C	132	PHE	O-C-N	-5.41	114.05	122.70
1	G	107	TRP	NE1-CE2-CZ2	-5.30	124.57	130.40
1	E	68	ASP	CB-CG-OD1	-5.29	113.54	118.30
1	E	107	TRP	CG-CD1-NE1	-5.28	104.82	110.10
1	C	116	ARG	NE-CZ-NH1	-5.25	117.67	120.30
1	H	37	ARG	NE-CZ-NH1	-5.24	117.68	120.30
1	F	42	ILE	O-C-N	-5.19	114.39	122.70
1	F	97	TYR	CB-CG-CD2	-5.12	117.93	121.00
1	C	107	TRP	CB-CG-CD1	-5.10	120.37	127.00
1	D	11	PRO	CA-N-CD	-5.08	104.39	111.50
1	E	122	ARG	CD-NE-CZ	-5.07	116.50	123.60
1	H	68	ASP	CB-CA-C	-5.05	100.31	110.40
1	B	107	TRP	CB-CG-CD1	-5.04	120.45	127.00

There are no chirality outliers.

All (620) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	10	LYS	Mainchain
1	A	100	LYS	Mainchain
1	A	103	ALA	Mainchain
1	A	104	GLU	Mainchain
1	A	105	LYS	Mainchain
1	A	108	PHE	Mainchain
1	A	11	PRO	Mainchain
1	A	110	GLY	Mainchain
1	A	114	ASN	Mainchain
1	A	115	GLY	Mainchain
1	A	116	ARG	Mainchain
1	A	12	LYS	Mainchain
1	A	120	GLY	Mainchain
1	A	122	ARG	Mainchain

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Mol	Chain	Res	Type	Group
1	A	123	THR	Mainchain
1	A	124	HIS	Mainchain
1	A	125	PHE	Sidechain,Mainchain
1	A	128	LYS	Mainchain
1	A	129	ALA	Mainchain
1	A	131	LEU	Mainchain
1	A	134	PRO	Mainchain
1	A	14	LEU	Mainchain
1	A	15	TYR	Sidechain
1	A	16	CYS	Mainchain
1	A	18	ASN	Mainchain
1	A	20	GLY	Mainchain
1	A	21	TYR	Mainchain
1	A	22	PHE	Mainchain
1	A	24	ARG	Sidechain,Mainchain
1	A	28	ASP	Mainchain
1	A	30	THR	Mainchain
1	A	31	VAL	Mainchain
1	A	33	GLY	Mainchain
1	A	35	LYS	Mainchain
1	A	36	ASP	Mainchain
1	A	38	SER	Mainchain
1	A	39	ASP	Sidechain,Mainchain
1	A	41	HIS	Mainchain
1	A	42	ILE	Mainchain
1	A	43	GLN	Mainchain
1	A	45	GLN	Mainchain
1	A	46	LEU	Mainchain
1	A	47	ALA	Mainchain
1	A	48	ALA	Mainchain
1	A	49	GLU	Mainchain
1	A	50	SER	Mainchain
1	A	51	ILE	Mainchain
1	A	53	GLU	Mainchain
1	A	54	VAL	Mainchain
1	A	55	TYR	Sidechain,Mainchain
1	A	56	ILE	Mainchain
1	A	58	SER	Mainchain
1	A	59	THR	Mainchain
1	A	60	GLU	Mainchain
1	A	61	THR	Mainchain
1	A	63	GLN	Mainchain

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Mol	Chain	Res	Type	Group
1	A	64	PHE	Mainchain
1	A	65	LEU	Mainchain
1	A	66	ALA	Mainchain
1	A	67	MET	Mainchain
1	A	69	THR	Mainchain
1	A	70	ASP	Mainchain
1	A	71	GLY	Mainchain
1	A	73	LEU	Mainchain
1	A	74	TYR	Sidechain,Mainchain
1	A	76	SER	Mainchain
1	A	77	GLN	Mainchain
1	A	79	PRO	Mainchain
1	A	80	ASN	Mainchain
1	A	82	GLU	Mainchain
1	A	84	LEU	Mainchain
1	A	86	LEU	Mainchain
1	A	87	GLU	Mainchain
1	A	88	ARG	Mainchain
1	A	90	GLU	Mainchain
1	A	91	GLU	Mainchain
1	A	92	ASN	Mainchain
1	A	93	HIS	Mainchain
1	A	95	ASN	Mainchain
1	A	96	THR	Mainchain
1	A	97	TYR	Sidechain,Mainchain
1	A	98	ILE	Mainchain
1	A	99	SER	Mainchain
1	B	102	HIS	Mainchain
1	B	103	ALA	Mainchain
1	B	105	LYS	Mainchain
1	B	108	PHE	Mainchain
1	B	11	PRO	Mainchain
1	B	110	GLY	Mainchain
1	B	113	LYS	Mainchain
1	B	114	ASN	Mainchain
1	B	117	SER	Mainchain
1	B	118	LYS	Mainchain
1	B	12	LYS	Mainchain
1	B	122	ARG	Sidechain
1	B	123	THR	Mainchain
1	B	13	LEU	Mainchain
1	B	14	LEU	Mainchain

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Mol	Chain	Res	Type	Group
1	B	16	CYS	Mainchain
1	B	17	SER	Mainchain
1	B	18	ASN	Mainchain
1	B	19	GLY	Mainchain
1	B	21	TYR	Mainchain
1	B	24	ARG	Sidechain,Mainchain
1	B	25	ILE	Mainchain
1	B	26	LEU	Mainchain
1	B	27	PRO	Mainchain
1	B	29	GLY	Mainchain
1	B	32	ASP	Mainchain
1	B	34	THR	Mainchain
1	B	36	ASP	Mainchain
1	B	37	ARG	Mainchain
1	B	38	SER	Mainchain
1	B	39	ASP	Sidechain,Mainchain
1	B	40	GLN	Mainchain
1	B	41	HIS	Mainchain
1	B	43	GLN	Mainchain
1	B	45	GLN	Mainchain
1	B	46	LEU	Mainchain
1	B	48	ALA	Mainchain
1	B	52	GLY	Mainchain
1	B	53	GLU	Mainchain
1	B	54	VAL	Mainchain
1	B	55	TYR	Sidechain,Mainchain
1	B	58	SER	Mainchain
1	B	59	THR	Mainchain
1	B	60	GLU	Mainchain
1	B	62	GLY	Mainchain
1	B	63	GLN	Mainchain
1	B	64	PHE	Mainchain
1	B	65	LEU	Mainchain
1	B	66	ALA	Mainchain
1	B	67	MET	Mainchain
1	B	68	ASP	Sidechain,Mainchain
1	B	69	THR	Mainchain
1	B	70	ASP	Sidechain
1	B	71	GLY	Mainchain
1	B	74	TYR	Sidechain,Mainchain
1	B	76	SER	Mainchain
1	B	79	PRO	Mainchain

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Mol	Chain	Res	Type	Group
1	B	80	ASN	Mainchain
1	B	83	CYS	Mainchain
1	B	84	LEU	Mainchain
1	B	85	PHE	Sidechain
1	B	86	LEU	Mainchain
1	B	87	GLU	Sidechain,Mainchain
1	B	88	ARG	Mainchain
1	B	89	LEU	Mainchain
1	B	92	ASN	Mainchain
1	B	93	HIS	Sidechain,Mainchain
1	B	95	ASN	Mainchain
1	B	97	TYR	Mainchain
1	B	99	SER	Mainchain
1	C	10	LYS	Mainchain
1	C	102	HIS	Mainchain
1	C	103	ALA	Mainchain
1	C	104	GLU	Mainchain
1	C	105	LYS	Mainchain
1	C	106	HIS	Mainchain
1	C	107	TRP	Mainchain
1	C	108	PHE	Sidechain
1	C	11	PRO	Mainchain
1	C	116	ARG	Sidechain,Mainchain
1	C	118	LYS	Mainchain
1	C	12	LYS	Mainchain
1	C	122	ARG	Mainchain
1	C	123	THR	Mainchain
1	C	125	PHE	Mainchain
1	C	129	ALA	Mainchain
1	C	13	LEU	Mainchain
1	C	131	LEU	Mainchain
1	C	132	PHE	Mainchain
1	C	134	PRO	Mainchain
1	C	135	LEU	Mainchain
1	C	15	TYR	Sidechain,Mainchain
1	C	16	CYS	Mainchain
1	C	20	GLY	Mainchain
1	C	21	TYR	Sidechain,Mainchain
1	C	23	LEU	Mainchain
1	C	25	ILE	Mainchain
1	C	27	PRO	Mainchain
1	C	29	GLY	Mainchain

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Mol	Chain	Res	Type	Group
1	C	30	THR	Mainchain
1	C	32	ASP	Sidechain,Mainchain
1	C	35	LYS	Mainchain
1	C	36	ASP	Sidechain,Mainchain
1	C	37	ARG	Mainchain
1	C	39	ASP	Sidechain,Mainchain
1	C	40	GLN	Mainchain
1	C	41	HIS	Mainchain
1	C	42	ILE	Mainchain
1	C	47	ALA	Mainchain
1	C	48	ALA	Mainchain
1	C	50	SER	Mainchain
1	C	52	GLY	Mainchain
1	C	53	GLU	Mainchain
1	C	55	TYR	Mainchain
1	C	56	ILE	Mainchain
1	C	57	LYS	Mainchain
1	C	58	SER	Mainchain
1	C	60	GLU	Mainchain
1	C	63	GLN	Mainchain
1	C	65	LEU	Mainchain
1	C	66	ALA	Mainchain
1	C	68	ASP	Sidechain
1	C	69	THR	Mainchain
1	C	71	GLY	Mainchain
1	C	74	TYR	Sidechain,Mainchain
1	C	77	GLN	Mainchain
1	C	79	PRO	Mainchain
1	C	80	ASN	Mainchain
1	C	82	GLU	Mainchain
1	C	84	LEU	Mainchain
1	C	85	PHE	Sidechain,Mainchain
1	C	87	GLU	Mainchain
1	C	92	ASN	Mainchain
1	C	93	HIS	Mainchain
1	C	94	TYR	Mainchain
1	C	95	ASN	Mainchain
1	C	97	TYR	Sidechain
1	C	99	SER	Mainchain
1	D	100	LYS	Mainchain
1	D	101	LYS	Mainchain
1	D	104	GLU	Mainchain

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Mol	Chain	Res	Type	Group
1	D	105	LYS	Mainchain
1	D	108	PHE	Mainchain
1	D	11	PRO	Mainchain
1	D	110	GLY	Mainchain
1	D	111	LEU	Mainchain
1	D	113	LYS	Mainchain
1	D	114	ASN	Mainchain
1	D	115	GLY	Mainchain
1	D	117	SER	Mainchain
1	D	12	LYS	Mainchain
1	D	122	ARG	Sidechain,Mainchain
1	D	123	THR	Mainchain
1	D	124	HIS	Sidechain
1	D	130	ILE	Mainchain
1	D	131	LEU	Mainchain
1	D	132	PHE	Mainchain
1	D	134	PRO	Mainchain
1	D	135	LEU	Mainchain
1	D	16	CYS	Mainchain
1	D	17	SER	Mainchain
1	D	21	TYR	Mainchain
1	D	24	ARG	Sidechain
1	D	26	LEU	Mainchain
1	D	27	PRO	Mainchain
1	D	28	ASP	Mainchain
1	D	30	THR	Mainchain
1	D	31	VAL	Mainchain
1	D	32	ASP	Mainchain
1	D	34	THR	Mainchain
1	D	35	LYS	Mainchain
1	D	36	ASP	Sidechain,Mainchain
1	D	37	ARG	Sidechain,Mainchain
1	D	39	ASP	Mainchain
1	D	41	HIS	Mainchain
1	D	42	ILE	Mainchain
1	D	43	GLN	Mainchain
1	D	44	LEU	Mainchain
1	D	46	LEU	Mainchain
1	D	48	ALA	Mainchain
1	D	49	GLU	Mainchain
1	D	50	SER	Mainchain
1	D	51	ILE	Mainchain

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Mol	Chain	Res	Type	Group
1	D	52	GLY	Mainchain
1	D	53	GLU	Sidechain,Mainchain
1	D	54	VAL	Mainchain
1	D	55	TYR	Mainchain
1	D	58	SER	Mainchain
1	D	59	THR	Mainchain
1	D	61	THR	Mainchain
1	D	64	PHE	Sidechain,Mainchain
1	D	65	LEU	Mainchain
1	D	70	ASP	Sidechain
1	D	72	LEU	Mainchain
1	D	75	GLY	Mainchain
1	D	76	SER	Mainchain
1	D	79	PRO	Mainchain
1	D	80	ASN	Mainchain
1	D	82	GLU	Mainchain
1	D	87	GLU	Sidechain,Mainchain
1	D	88	ARG	Mainchain
1	D	90	GLU	Mainchain
1	D	92	ASN	Mainchain
1	D	95	ASN	Mainchain
1	D	96	THR	Mainchain
1	D	97	TYR	Sidechain,Mainchain
1	D	99	SER	Mainchain
1	E	10	LYS	Mainchain
1	E	102	HIS	Mainchain
1	E	103	ALA	Mainchain
1	E	104	GLU	Sidechain,Mainchain
1	E	106	HIS	Mainchain
1	E	107	TRP	Mainchain
1	E	108	PHE	Mainchain
1	E	11	PRO	Mainchain
1	E	110	GLY	Mainchain
1	E	111	LEU	Mainchain
1	E	113	LYS	Mainchain
1	E	114	ASN	Mainchain
1	E	115	GLY	Mainchain
1	E	117	SER	Mainchain
1	E	118	LYS	Mainchain
1	E	12	LYS	Mainchain
1	E	125	PHE	Mainchain
1	E	129	ALA	Mainchain

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Mol	Chain	Res	Type	Group
1	E	13	LEU	Mainchain
1	E	130	ILE	Mainchain
1	E	131	LEU	Mainchain
1	E	135	LEU	Mainchain
1	E	15	TYR	Mainchain
1	E	20	GLY	Mainchain
1	E	22	PHE	Mainchain
1	E	23	LEU	Mainchain
1	E	25	ILE	Mainchain
1	E	26	LEU	Mainchain
1	E	27	PRO	Mainchain
1	E	28	ASP	Mainchain
1	E	31	VAL	Mainchain
1	E	32	ASP	Sidechain
1	E	36	ASP	Mainchain
1	E	37	ARG	Sidechain,Mainchain
1	E	39	ASP	Mainchain
1	E	40	GLN	Mainchain
1	E	41	HIS	Mainchain
1	E	43	GLN	Mainchain
1	E	45	GLN	Mainchain
1	E	46	LEU	Mainchain
1	E	47	ALA	Mainchain
1	E	49	GLU	Mainchain
1	E	52	GLY	Mainchain
1	E	53	GLU	Mainchain
1	E	54	VAL	Mainchain
1	E	55	TYR	Mainchain
1	E	58	SER	Mainchain
1	E	59	THR	Mainchain
1	E	63	GLN	Mainchain
1	E	64	PHE	Mainchain
1	E	65	LEU	Mainchain
1	E	67	MET	Mainchain
1	E	68	ASP	Mainchain
1	E	69	THR	Mainchain
1	E	71	GLY	Mainchain
1	E	72	LEU	Mainchain
1	E	74	TYR	Mainchain
1	E	77	GLN	Mainchain
1	E	79	PRO	Mainchain
1	E	80	ASN	Mainchain

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Mol	Chain	Res	Type	Group
1	E	82	GLU	Mainchain
1	E	83	CYS	Mainchain
1	E	85	PHE	Sidechain
1	E	86	LEU	Mainchain
1	E	87	GLU	Sidechain
1	E	88	ARG	Mainchain
1	E	89	LEU	Mainchain
1	E	91	GLU	Mainchain
1	E	92	ASN	Mainchain
1	E	93	HIS	Sidechain
1	E	96	THR	Mainchain
1	E	97	TYR	Sidechain
1	F	10	LYS	Mainchain
1	F	100	LYS	Mainchain
1	F	101	LYS	Mainchain
1	F	103	ALA	Mainchain
1	F	104	GLU	Mainchain
1	F	107	TRP	Mainchain
1	F	108	PHE	Mainchain
1	F	110	GLY	Mainchain
1	F	112	LYS	Mainchain
1	F	115	GLY	Mainchain
1	F	117	SER	Mainchain
1	F	118	LYS	Mainchain
1	F	12	LYS	Mainchain
1	F	121	PRO	Mainchain
1	F	122	ARG	Sidechain
1	F	123	THR	Mainchain
1	F	125	PHE	Sidechain
1	F	128	LYS	Mainchain
1	F	132	PHE	Mainchain
1	F	14	LEU	Mainchain
1	F	15	TYR	Sidechain
1	F	20	GLY	Mainchain
1	F	21	TYR	Mainchain
1	F	22	PHE	Sidechain,Mainchain
1	F	23	LEU	Mainchain
1	F	24	ARG	Mainchain
1	F	25	ILE	Mainchain
1	F	30	THR	Mainchain
1	F	31	VAL	Mainchain
1	F	32	ASP	Sidechain

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Mol	Chain	Res	Type	Group
1	F	33	GLY	Mainchain
1	F	34	THR	Mainchain
1	F	36	ASP	Mainchain
1	F	37	ARG	Mainchain
1	F	40	GLN	Mainchain
1	F	42	ILE	Mainchain
1	F	43	GLN	Mainchain
1	F	44	LEU	Mainchain
1	F	46	LEU	Mainchain
1	F	47	ALA	Mainchain
1	F	48	ALA	Mainchain
1	F	49	GLU	Mainchain
1	F	50	SER	Mainchain
1	F	51	ILE	Mainchain
1	F	52	GLY	Mainchain
1	F	53	GLU	Mainchain
1	F	54	VAL	Mainchain
1	F	55	TYR	Sidechain,Mainchain
1	F	56	ILE	Mainchain
1	F	60	GLU	Mainchain
1	F	63	GLN	Mainchain
1	F	65	LEU	Mainchain
1	F	66	ALA	Mainchain
1	F	68	ASP	Mainchain
1	F	70	ASP	Sidechain
1	F	73	LEU	Mainchain
1	F	74	TYR	Mainchain
1	F	75	GLY	Mainchain
1	F	76	SER	Mainchain
1	F	78	THR	Mainchain
1	F	79	PRO	Mainchain
1	F	82	GLU	Mainchain
1	F	83	CYS	Mainchain
1	F	85	PHE	Sidechain,Mainchain
1	F	86	LEU	Mainchain
1	F	87	GLU	Sidechain
1	F	88	ARG	Mainchain
1	F	90	GLU	Mainchain
1	F	92	ASN	Mainchain
1	F	94	TYR	Mainchain
1	F	95	ASN	Mainchain
1	F	96	THR	Mainchain

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Mol	Chain	Res	Type	Group
1	F	97	TYR	Sidechain,Mainchain
1	F	98	ILE	Mainchain
1	G	100	LYS	Mainchain
1	G	105	LYS	Mainchain
1	G	106	HIS	Mainchain
1	G	108	PHE	Mainchain
1	G	11	PRO	Mainchain
1	G	112	LYS	Mainchain
1	G	113	LYS	Mainchain
1	G	114	ASN	Mainchain
1	G	115	GLY	Mainchain
1	G	119	LEU	Mainchain
1	G	12	LYS	Mainchain
1	G	121	PRO	Mainchain
1	G	123	THR	Mainchain
1	G	128	LYS	Mainchain
1	G	13	LEU	Mainchain
1	G	132	PHE	Mainchain
1	G	135	LEU	Mainchain
1	G	15	TYR	Sidechain
1	G	16	CYS	Mainchain
1	G	21	TYR	Mainchain
1	G	22	PHE	Mainchain
1	G	24	ARG	Sidechain,Mainchain
1	G	25	ILE	Mainchain
1	G	26	LEU	Mainchain
1	G	29	GLY	Mainchain
1	G	32	ASP	Mainchain
1	G	33	GLY	Mainchain
1	G	36	ASP	Mainchain
1	G	38	SER	Mainchain
1	G	39	ASP	Mainchain
1	G	40	GLN	Mainchain
1	G	45	GLN	Mainchain
1	G	47	ALA	Mainchain
1	G	52	GLY	Mainchain
1	G	53	GLU	Sidechain,Mainchain
1	G	54	VAL	Mainchain
1	G	55	TYR	Mainchain
1	G	57	LYS	Mainchain
1	G	58	SER	Mainchain
1	G	59	THR	Mainchain

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Mol	Chain	Res	Type	Group
1	G	60	GLU	Mainchain
1	G	64	PHE	Sidechain,Mainchain
1	G	67	MET	Mainchain
1	G	69	THR	Mainchain
1	G	70	ASP	Mainchain
1	G	71	GLY	Mainchain
1	G	72	LEU	Mainchain
1	G	74	TYR	Mainchain
1	G	75	GLY	Mainchain
1	G	77	GLN	Mainchain
1	G	79	PRO	Mainchain
1	G	82	GLU	Mainchain
1	G	83	CYS	Mainchain
1	G	85	PHE	Mainchain
1	G	86	LEU	Mainchain
1	G	87	GLU	Sidechain,Mainchain
1	G	88	ARG	Mainchain
1	G	92	ASN	Mainchain
1	G	93	HIS	Mainchain
1	G	95	ASN	Mainchain
1	G	97	TYR	Mainchain
1	G	98	ILE	Mainchain
1	G	99	SER	Mainchain
1	H	100	LYS	Mainchain
1	H	102	HIS	Mainchain
1	H	103	ALA	Mainchain
1	H	104	GLU	Mainchain
1	H	105	LYS	Mainchain
1	H	108	PHE	Mainchain
1	H	109	VAL	Mainchain
1	H	111	LEU	Mainchain
1	H	113	LYS	Mainchain
1	H	115	GLY	Mainchain
1	H	116	ARG	Mainchain
1	H	118	LYS	Mainchain
1	H	12	LYS	Mainchain
1	H	123	THR	Mainchain
1	H	124	HIS	Sidechain,Mainchain
1	H	125	PHE	Sidechain,Mainchain
1	H	129	ALA	Mainchain
1	H	13	LEU	Mainchain
1	H	130	ILE	Mainchain

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Mol	Chain	Res	Type	Group
1	H	131	LEU	Mainchain
1	H	132	PHE	Mainchain
1	H	134	PRO	Mainchain
1	H	17	SER	Mainchain
1	H	18	ASN	Mainchain
1	H	20	GLY	Mainchain
1	H	21	TYR	Sidechain,Mainchain
1	H	27	PRO	Mainchain
1	H	28	ASP	Sidechain
1	H	30	THR	Mainchain
1	H	32	ASP	Mainchain
1	H	35	LYS	Mainchain
1	H	37	ARG	Mainchain
1	H	38	SER	Mainchain
1	H	39	ASP	Sidechain,Mainchain
1	H	40	GLN	Mainchain
1	H	41	HIS	Mainchain
1	H	43	GLN	Mainchain
1	H	44	LEU	Mainchain
1	H	45	GLN	Mainchain
1	H	47	ALA	Mainchain
1	H	48	ALA	Mainchain
1	H	50	SER	Mainchain
1	H	51	ILE	Mainchain
1	H	52	GLY	Mainchain
1	H	53	GLU	Mainchain
1	H	55	TYR	Sidechain,Mainchain
1	H	56	ILE	Mainchain
1	H	57	LYS	Mainchain
1	H	58	SER	Mainchain
1	H	59	THR	Mainchain
1	H	60	GLU	Mainchain
1	H	61	THR	Mainchain
1	H	62	GLY	Mainchain
1	H	63	GLN	Mainchain
1	H	68	ASP	Mainchain
1	H	70	ASP	Mainchain
1	H	72	LEU	Mainchain
1	H	74	TYR	Sidechain,Mainchain
1	H	76	SER	Mainchain
1	H	77	GLN	Mainchain
1	H	79	PRO	Mainchain

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Mol	Chain	Res	Type	Group
1	H	80	ASN	Mainchain
1	H	81	GLU	Mainchain
1	H	84	LEU	Mainchain
1	H	85	PHE	Sidechain,Mainchain
1	H	86	LEU	Mainchain
1	H	88	ARG	Mainchain
1	H	89	LEU	Mainchain
1	H	90	GLU	Mainchain
1	H	91	GLU	Mainchain
1	H	92	ASN	Mainchain
1	H	93	HIS	Mainchain
1	H	94	TYR	Mainchain
1	H	96	THR	Mainchain
1	H	97	TYR	Mainchain
1	H	99	SER	Mainchain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	983	0	944	109	0
1	B	983	0	944	96	0
1	C	983	0	944	112	0
1	D	983	0	944	96	0
1	E	983	0	944	86	0
1	F	983	0	944	84	0
1	G	983	0	944	126	0
1	H	983	0	944	121	0
2	A	55	0	14	29	0
2	B	55	0	15	11	0
2	C	55	0	14	16	0
2	D	55	0	16	24	0
2	E	55	0	14	26	0
2	F	55	0	14	14	0
2	G	55	0	14	35	0
2	H	55	0	14	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	8304	0	7667	934	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 58.

All (934) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:E:141:SCR:H14	2:E:141:SCR:S14	1.82	1.20
1:A:114:ASN:HD21	1:A:116:ARG:NH1	1.43	1.16
1:G:24:ARG:HH11	1:G:26:LEU:HD11	1.01	1.14
1:H:27:PRO:HB3	1:H:61:THR:HG21	1.31	1.12
1:D:86:LEU:HD21	1:D:100:LYS:HG3	1.30	1.11
1:F:24:ARG:HD2	1:F:26:LEU:HD11	1.30	1.11
1:B:26:LEU:HB2	1:B:30:THR:HG22	1.22	1.11
1:D:98:ILE:HD11	1:D:106:HIS:HA	1.32	1.11
1:E:26:LEU:HD23	1:E:30:THR:CG2	1.85	1.06
2:G:141:SCR:H15	2:G:141:SCR:O72	1.25	1.05
1:E:95:ASN:N	1:E:95:ASN:HD22	1.51	1.05
2:A:141:SCR:C5	2:A:141:SCR:H162	1.87	1.04
2:A:141:SCR:C6	2:A:141:SCR:H162	1.87	1.04
1:E:26:LEU:HD23	1:E:30:THR:HG22	1.06	1.04
1:B:24:ARG:HD3	1:B:26:LEU:HD13	1.36	1.02
1:H:12:LYS:HB3	1:H:135:LEU:O	1.60	1.02
1:A:24:ARG:HD2	1:A:26:LEU:HD11	1.42	1.00
1:G:102:HIS:HB3	1:G:105:LYS:HG3	1.39	1.00
1:G:112:LYS:CE	2:G:141:SCR:H14	1.92	0.99
1:F:112:LYS:HE2	2:F:141:SCR:S12	2.03	0.99
1:G:24:ARG:NH1	1:G:26:LEU:HD11	1.78	0.99
1:G:26:LEU:HD23	1:G:30:THR:CG2	1.92	0.99
2:G:141:SCR:H62	2:G:141:SCR:O44	1.63	0.98
1:C:89:LEU:HD13	1:C:90:GLU:N	1.78	0.98
1:D:123:THR:O	1:D:124:HIS:HB3	1.60	0.98
1:E:26:LEU:CD2	1:E:30:THR:HG22	1.93	0.98
2:G:141:SCR:H112	2:G:141:SCR:O72	1.63	0.97
2:E:141:SCR:H14	2:E:141:SCR:O52	1.65	0.97
1:F:92:ASN:O	1:F:93:HIS:HB2	1.64	0.96
1:H:53:GLU:HG2	1:H:84:LEU:HD23	1.47	0.94
2:G:141:SCR:S6	2:G:141:SCR:H161	2.07	0.94
1:G:26:LEU:HD23	1:G:30:THR:HG22	1.46	0.94
1:A:90:GLU:HA	1:A:90:GLU:OE1	1.68	0.94
1:C:55:TYR:CZ	1:C:84:LEU:HD21	2.05	0.91
1:B:13:LEU:HB2	1:B:135:LEU:O	1.69	0.91

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:141:SCR:O54	2:H:141:SCR:H62	1.71	0.91
1:G:112:LYS:HE2	2:G:141:SCR:H14	1.48	0.91
1:A:89:LEU:HD23	1:A:90:GLU:N	1.85	0.91
1:D:24:ARG:HD3	1:D:26:LEU:HD13	1.50	0.91
1:C:108:PHE:CE1	1:C:130:ILE:HD11	2.06	0.91
1:A:11:PRO:O	1:A:12:LYS:HD3	1.71	0.91
1:C:65:LEU:HD21	1:C:73:LEU:HD23	1.50	0.90
1:A:46:LEU:HD23	1:A:56:ILE:HG12	1.54	0.90
2:A:141:SCR:H162	2:A:141:SCR:H61	1.54	0.90
1:B:16:CYS:HB2	1:B:132:PHE:CE1	2.07	0.90
1:A:26:LEU:N	1:A:26:LEU:HD12	1.88	0.89
2:G:141:SCR:O6	2:G:141:SCR:H161	1.71	0.89
1:C:24:ARG:HD3	1:C:26:LEU:HD13	1.53	0.89
1:B:28:ASP:OD1	1:B:30:THR:HB	1.74	0.88
1:H:12:LYS:CB	1:H:135:LEU:O	2.22	0.88
1:B:112:LYS:HE3	2:B:141:SCR:O92	1.73	0.88
1:A:114:ASN:ND2	1:A:116:ARG:NH1	2.22	0.87
1:G:80:ASN:O	1:G:83:CYS:HB2	1.74	0.87
2:G:141:SCR:C15	2:G:141:SCR:O72	2.16	0.87
1:E:86:LEU:HD21	1:E:100:LYS:CD	2.05	0.86
2:C:141:SCR:H111	2:C:141:SCR:O2	1.75	0.86
1:G:24:ARG:HD3	1:G:26:LEU:HD13	1.55	0.86
1:D:98:ILE:HD11	1:D:106:HIS:CA	2.05	0.86
2:A:141:SCR:C16	2:A:141:SCR:H61	2.06	0.85
1:A:46:LEU:HB3	1:A:56:ILE:HG13	1.58	0.85
1:G:124:HIS:C	1:G:124:HIS:CD2	2.49	0.85
1:D:48:ALA:HA	1:D:54:VAL:HG12	1.57	0.85
1:B:56:ILE:HD12	1:B:65:LEU:HD12	1.58	0.85
1:H:112:LYS:CE	2:H:141:SCR:H61	2.07	0.85
1:A:102:HIS:HA	1:A:104:GLU:OE1	1.76	0.84
1:D:96:THR:HG23	1:D:130:ILE:HG23	1.58	0.84
1:G:24:ARG:HH11	1:G:26:LEU:CD1	1.87	0.84
1:A:89:LEU:HD23	1:A:90:GLU:H	1.42	0.84
1:F:53:GLU:HB3	1:F:84:LEU:HD22	1.60	0.83
1:F:86:LEU:HD21	1:F:100:LYS:HD2	1.61	0.83
1:G:124:HIS:O	1:G:124:HIS:HD2	1.61	0.83
1:G:112:LYS:HE2	2:G:141:SCR:C16	2.09	0.82
1:D:112:LYS:HZ1	2:D:141:SCR:H61	1.44	0.82
1:G:86:LEU:HD21	1:G:100:LYS:HG3	1.59	0.82
2:A:141:SCR:H14	2:A:141:SCR:O52	1.79	0.82
1:D:108:PHE:O	1:D:123:THR:HG21	1.80	0.82
1:D:112:LYS:NZ	2:D:141:SCR:H61	1.95	0.82

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:57:LYS:HG3	1:H:58:SER:N	1.94	0.82
1:D:77:GLN:O	1:D:77:GLN:HG3	1.78	0.82
1:B:102:HIS:HB3	1:B:105:LYS:HG3	1.60	0.82
1:B:89:LEU:HD12	1:B:90:GLU:N	1.94	0.82
2:C:141:SCR:S2	2:C:141:SCR:H111	2.19	0.82
1:A:26:LEU:N	1:A:26:LEU:CD1	2.43	0.81
1:H:111:LEU:HD23	1:H:117:SER:HA	1.61	0.81
2:A:141:SCR:S14	2:A:141:SCR:H61	2.19	0.81
1:F:89:LEU:HD21	1:F:93:HIS:ND1	1.94	0.81
1:C:122:ARG:CD	2:C:141:SCR:O83	2.28	0.81
1:C:104:GLU:OE2	1:C:105:LYS:HD3	1.79	0.81
1:B:64:PHE:CZ	1:B:79:PRO:HG3	2.15	0.80
2:A:141:SCR:C6	2:A:141:SCR:C16	2.59	0.80
1:B:94:TYR:CE1	1:B:133:LEU:HB2	2.16	0.80
1:C:60:GLU:N	1:C:60:GLU:OE1	2.15	0.80
1:F:112:LYS:HE2	2:F:141:SCR:O94	1.80	0.80
1:C:24:ARG:HD2	1:C:32:ASP:OD1	1.81	0.80
1:B:39:ASP:O	1:B:42:ILE:HG12	1.80	0.80
1:E:39:ASP:OD1	1:E:41:HIS:HB2	1.81	0.80
1:D:112:LYS:HE2	2:D:141:SCR:S12	2.23	0.79
1:D:86:LEU:CD2	1:D:100:LYS:HG3	2.10	0.79
1:F:124:HIS:C	1:F:130:ILE:HD11	2.03	0.79
2:E:141:SCR:S14	2:E:141:SCR:C14	2.67	0.79
2:G:141:SCR:O63	2:G:141:SCR:H161	1.83	0.79
1:B:72:LEU:HD23	1:B:119:LEU:HD21	1.63	0.79
1:F:49:GLU:HG3	1:F:84:LEU:HD21	1.63	0.79
1:G:112:LYS:HE2	2:G:141:SCR:C14	2.13	0.78
1:D:37:ARG:O	1:D:37:ARG:HG2	1.84	0.78
1:B:26:LEU:CB	1:B:30:THR:HG22	2.11	0.78
1:H:57:LYS:HG3	1:H:58:SER:H	1.49	0.78
1:C:131:LEU:H	1:C:131:LEU:HD22	1.49	0.78
1:A:84:LEU:HB2	1:A:100:LYS:CE	2.13	0.77
1:C:12:LYS:HA	1:C:136:PRO:HA	1.65	0.77
2:E:141:SCR:S2	2:E:141:SCR:H112	2.23	0.77
1:A:124:HIS:CE1	1:A:127:GLN:NE2	2.53	0.77
1:H:90:GLU:O	1:H:91:GLU:CB	2.30	0.77
1:G:105:LYS:O	1:G:106:HIS:HB2	1.84	0.77
2:F:141:SCR:O54	2:F:141:SCR:O6	2.02	0.77
1:G:39:ASP:O	1:G:42:ILE:HG12	1.84	0.77
1:D:14:LEU:HD22	1:D:44:LEU:HD23	1.67	0.77
1:D:130:ILE:HG22	1:D:131:LEU:HD13	1.66	0.77
1:G:89:LEU:HD22	1:G:90:GLU:H	1.49	0.77

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:24:ARG:HD2	1:F:26:LEU:CD1	2.13	0.77
1:A:24:ARG:CD	1:A:26:LEU:HD11	2.14	0.77
2:A:141:SCR:H14	2:A:141:SCR:S14	2.25	0.76
1:A:90:GLU:CA	1:A:90:GLU:OE1	2.31	0.76
1:C:122:ARG:HD2	2:C:141:SCR:O83	1.84	0.76
1:A:105:LYS:O	1:A:106:HIS:HB2	1.84	0.76
1:G:124:HIS:CD2	1:G:125:PHE:C	2.59	0.76
1:G:124:HIS:CD2	1:G:124:HIS:O	2.39	0.76
2:E:141:SCR:H14	2:E:141:SCR:O54	1.84	0.76
1:C:12:LYS:HB3	1:C:135:LEU:O	1.86	0.76
1:E:84:LEU:O	1:E:100:LYS:HB2	1.86	0.76
1:E:90:GLU:HB2	1:E:94:TYR:O	1.86	0.75
2:C:141:SCR:O22	2:C:141:SCR:O84	2.04	0.75
1:B:94:TYR:HE1	1:B:133:LEU:HB2	1.51	0.75
1:G:61:THR:OG1	1:G:63:GLN:NE2	2.18	0.75
1:C:54:VAL:O	1:C:84:LEU:HD13	1.86	0.75
1:F:89:LEU:HD21	1:F:93:HIS:CE1	2.22	0.75
1:D:112:LYS:HG2	2:D:141:SCR:O92	1.85	0.75
2:F:141:SCR:O52	2:F:141:SCR:O63	2.05	0.75
1:G:65:LEU:O	1:G:66:ALA:HB2	1.85	0.75
1:F:46:LEU:HD23	1:F:56:ILE:HG13	1.67	0.75
1:G:24:ARG:HD3	1:G:26:LEU:CD1	2.17	0.75
1:F:89:LEU:O	1:F:90:GLU:HG2	1.86	0.75
1:H:53:GLU:HG3	1:H:86:LEU:HD23	1.69	0.74
1:H:123:THR:O	1:H:124:HIS:HB3	1.88	0.74
1:C:108:PHE:CZ	1:C:130:ILE:HD11	2.23	0.74
1:F:26:LEU:HD13	1:F:26:LEU:N	2.03	0.74
1:H:112:LYS:HE3	2:H:141:SCR:H61	1.67	0.73
1:E:112:LYS:HE2	2:E:141:SCR:O53	1.88	0.73
2:E:141:SCR:O24	2:E:141:SCR:H112	1.89	0.73
1:A:112:LYS:HE3	2:A:141:SCR:O73	1.88	0.73
1:G:26:LEU:HB2	1:G:30:THR:HG22	1.71	0.73
1:D:98:ILE:CD1	1:D:106:HIS:HA	2.15	0.73
1:G:87:GLU:HG3	1:G:87:GLU:O	1.87	0.73
2:F:141:SCR:O84	2:F:141:SCR:H3	1.89	0.73
1:D:87:GLU:HG3	1:D:97:TYR:CE2	2.23	0.73
1:C:37:ARG:O	1:C:37:ARG:HG2	1.88	0.72
1:A:23:LEU:HD23	1:A:44:LEU:HD21	1.70	0.72
1:D:96:THR:CG2	1:D:130:ILE:HG23	2.18	0.72
1:B:56:ILE:CD1	1:B:65:LEU:HD12	2.17	0.72
1:D:19:GLY:HA3	1:D:21:TYR:CE2	2.24	0.72
1:C:108:PHE:CD1	1:C:130:ILE:HD11	2.25	0.71

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:112:LYS:CE	2:G:141:SCR:H162	2.20	0.71
1:G:124:HIS:C	1:G:124:HIS:HD2	1.90	0.71
1:A:118:LYS:HE2	2:A:141:SCR:O92	1.91	0.71
1:B:86:LEU:HD21	1:B:100:LYS:CG	2.20	0.71
1:B:26:LEU:HB2	1:B:30:THR:CG2	2.13	0.71
1:F:111:LEU:HD23	1:F:117:SER:HA	1.73	0.71
1:G:16:CYS:HB2	1:G:132:PHE:CE1	2.26	0.71
1:E:105:LYS:O	1:E:106:HIS:HB2	1.89	0.71
1:G:112:LYS:HD3	2:G:141:SCR:O63	1.91	0.71
1:C:55:TYR:OH	1:C:84:LEU:HD21	1.90	0.71
1:E:89:LEU:HG	1:E:90:GLU:H	1.57	0.70
1:C:46:LEU:HD22	1:C:56:ILE:HG12	1.73	0.70
2:H:141:SCR:O24	2:H:141:SCR:H111	1.91	0.70
1:C:104:GLU:H	1:C:104:GLU:CD	1.95	0.70
1:C:89:LEU:O	1:C:90:GLU:HG2	1.92	0.70
1:C:16:CYS:HB2	1:C:132:PHE:CE2	2.27	0.70
2:H:141:SCR:H111	2:H:141:SCR:S2	2.32	0.70
2:B:141:SCR:O93	2:B:141:SCR:O81	2.09	0.70
2:C:141:SCR:H111	2:C:141:SCR:O24	1.90	0.70
1:H:79:PRO:O	1:H:80:ASN:HB3	1.92	0.70
2:H:141:SCR:O54	2:H:141:SCR:C6	2.39	0.70
1:B:92:ASN:O	1:B:93:HIS:HB2	1.91	0.70
1:G:89:LEU:HD22	1:G:90:GLU:N	2.06	0.69
2:H:141:SCR:S14	2:H:141:SCR:H62	2.31	0.69
1:C:122:ARG:HD3	2:C:141:SCR:O83	1.90	0.69
1:E:64:PHE:CE1	1:E:79:PRO:HD3	2.26	0.69
1:D:15:TYR:HB2	1:D:135:LEU:CD1	2.21	0.69
2:G:141:SCR:O91	2:G:141:SCR:H3	1.91	0.69
1:E:55:TYR:OH	1:E:84:LEU:HD22	1.93	0.69
1:A:84:LEU:CB	1:A:100:LYS:CE	2.70	0.69
1:G:49:GLU:HG3	1:G:55:TYR:CE2	2.26	0.69
1:F:39:ASP:HB3	1:F:42:ILE:HG23	1.74	0.69
1:A:114:ASN:HD21	1:A:116:ARG:HH12	1.40	0.69
1:A:85:PHE:HA	1:A:98:ILE:O	1.93	0.69
1:F:26:LEU:H	1:F:26:LEU:HD13	1.56	0.68
1:F:24:ARG:HG3	1:F:34:THR:HG21	1.75	0.68
1:D:15:TYR:HB2	1:D:135:LEU:HD12	1.76	0.68
1:E:43:GLN:HE22	1:E:60:GLU:HG3	1.57	0.68
1:A:112:LYS:NZ	2:A:141:SCR:O92	2.27	0.68
2:G:141:SCR:H112	2:G:141:SCR:S13	2.34	0.68
1:C:127:GLN:O	1:C:130:ILE:HG22	1.93	0.68
1:B:72:LEU:HD23	1:B:119:LEU:CD2	2.23	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:27:PRO:HB3	1:H:61:THR:CG2	2.19	0.67
2:F:141:SCR:O81	2:F:141:SCR:O93	2.11	0.67
1:F:47:ALA:O	1:F:55:TYR:HD2	1.76	0.67
1:G:47:ALA:C	1:G:54:VAL:HG12	2.14	0.67
2:E:141:SCR:C14	2:E:141:SCR:O52	2.40	0.67
1:B:64:PHE:CE2	1:B:79:PRO:HG3	2.28	0.67
1:C:131:LEU:O	1:C:131:LEU:HD23	1.94	0.67
1:A:84:LEU:HB2	1:A:100:LYS:HD2	1.76	0.67
1:G:124:HIS:HD2	1:G:125:PHE:C	1.97	0.67
1:C:73:LEU:HD21	1:C:109:VAL:HG13	1.76	0.67
1:F:26:LEU:CD1	1:F:26:LEU:N	2.57	0.67
1:G:112:LYS:HE2	2:G:141:SCR:H162	1.74	0.67
1:G:26:LEU:CD2	1:G:30:THR:HG22	2.24	0.67
1:E:26:LEU:HB2	1:E:28:ASP:OD1	1.95	0.67
1:A:111:LEU:HD23	1:A:117:SER:HA	1.75	0.67
1:E:95:ASN:ND2	1:E:95:ASN:N	2.25	0.67
1:A:80:ASN:HD22	1:A:82:GLU:H	1.41	0.67
1:A:46:LEU:HA	1:A:55:TYR:O	1.95	0.67
2:E:141:SCR:O22	2:E:141:SCR:O94	2.13	0.66
1:C:108:PHE:HB2	1:C:123:THR:HB	1.77	0.66
1:C:65:LEU:HD21	1:C:73:LEU:CD2	2.23	0.66
1:F:124:HIS:O	1:F:130:ILE:HD11	1.94	0.66
1:H:105:LYS:O	1:H:106:HIS:HB2	1.94	0.66
2:E:141:SCR:H13	2:E:141:SCR:O74	1.94	0.66
1:C:134:PRO:C	1:C:135:LEU:HD23	2.15	0.66
1:D:87:GLU:O	1:D:88:ARG:CB	2.43	0.66
1:B:49:GLU:OE1	1:B:53:GLU:HB3	1.95	0.66
1:C:128:LYS:HB2	1:C:128:LYS:NZ	2.08	0.66
1:C:46:LEU:HA	1:C:55:TYR:O	1.95	0.66
1:G:66:ALA:HB1	1:G:74:TYR:CE1	2.30	0.66
1:G:49:GLU:HG3	1:G:55:TYR:HE2	1.61	0.66
1:F:105:LYS:O	1:F:106:HIS:HB2	1.94	0.66
1:A:94:TYR:CD2	1:A:133:LEU:HB2	2.31	0.66
1:D:98:ILE:HD12	1:D:103:ALA:HA	1.78	0.66
1:A:54:VAL:HG22	1:A:85:PHE:O	1.96	0.66
1:C:98:ILE:HD11	1:C:108:PHE:CE1	2.30	0.66
1:C:12:LYS:HA	1:C:135:LEU:O	1.96	0.66
2:G:141:SCR:S4	2:G:141:SCR:H62	2.36	0.65
1:G:57:LYS:HG3	1:G:57:LYS:O	1.95	0.65
1:G:112:LYS:CE	2:G:141:SCR:C14	2.70	0.65
2:G:141:SCR:S12	2:G:141:SCR:H3	2.36	0.65
1:B:24:ARG:HD3	1:B:26:LEU:CD1	2.20	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:79:PRO:O	1:B:80:ASN:HB3	1.95	0.65
1:B:86:LEU:HD21	1:B:100:LYS:HG3	1.79	0.65
1:B:24:ARG:HD2	1:B:32:ASP:OD1	1.96	0.65
1:G:102:HIS:HB3	1:G:105:LYS:CG	2.20	0.65
2:C:141:SCR:H62	2:C:141:SCR:H161	1.79	0.65
1:H:104:GLU:OE1	1:H:105:LYS:N	2.29	0.65
1:F:12:LYS:HD3	1:F:135:LEU:O	1.96	0.65
1:C:64:PHE:CZ	1:C:79:PRO:HG3	2.31	0.65
1:F:36:ASP:C	1:F:38:SER:H	1.98	0.65
2:D:141:SCR:O5	2:D:141:SCR:H161	1.96	0.65
2:A:141:SCR:O44	2:A:141:SCR:O34	2.15	0.65
1:E:68:ASP:HB3	1:E:70:ASP:H	1.60	0.65
1:F:68:ASP:OD1	1:F:72:LEU:HB3	1.96	0.65
1:A:112:LYS:HB3	1:A:112:LYS:NZ	2.11	0.65
1:E:80:ASN:ND2	1:E:83:CYS:SG	2.70	0.65
1:C:23:LEU:HD21	1:C:65:LEU:HD11	1.78	0.64
1:G:124:HIS:CD2	1:G:125:PHE:O	2.50	0.64
2:C:141:SCR:H62	2:C:141:SCR:O54	1.97	0.64
1:H:24:ARG:NE	1:H:26:LEU:HD11	2.13	0.64
1:C:47:ALA:HB3	1:C:55:TYR:HB2	1.78	0.64
1:D:24:ARG:HD3	1:D:26:LEU:CD1	2.26	0.64
2:A:141:SCR:H5	2:A:141:SCR:H162	1.79	0.64
1:B:68:ASP:HB3	1:B:72:LEU:H	1.62	0.64
1:H:46:LEU:HD12	1:H:46:LEU:H	1.62	0.64
1:C:46:LEU:HD23	1:C:46:LEU:N	2.12	0.64
1:F:37:ARG:HH21	1:F:135:LEU:HD13	1.63	0.64
1:D:64:PHE:HE1	1:D:78:THR:HA	1.62	0.64
1:G:65:LEU:O	1:G:66:ALA:CB	2.45	0.63
1:C:55:TYR:CZ	1:C:84:LEU:CD2	2.82	0.63
1:G:64:PHE:CE2	1:G:79:PRO:N	2.66	0.63
1:A:31:VAL:O	1:A:117:SER:HB3	1.97	0.63
1:H:24:ARG:HD3	1:H:26:LEU:HD13	1.79	0.63
1:E:104:GLU:OE2	1:E:105:LYS:N	2.30	0.63
2:B:141:SCR:O1	2:B:141:SCR:O93	2.17	0.63
1:H:57:LYS:CG	1:H:58:SER:N	2.61	0.63
2:E:141:SCR:H112	2:E:141:SCR:O2	1.99	0.63
1:D:55:TYR:CE1	1:D:79:PRO:HB3	2.33	0.63
1:B:87:GLU:HG3	1:B:88:ARG:N	2.13	0.63
1:D:114:ASN:OD1	1:D:116:ARG:HB2	1.98	0.63
2:A:141:SCR:O51	2:A:141:SCR:H61	1.97	0.62
1:H:129:ALA:HB3	1:H:130:ILE:HD13	1.81	0.62
1:H:24:ARG:N	1:H:32:ASP:O	2.28	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:89:LEU:HD11	1:B:93:HIS:HA	1.80	0.62
1:E:58:SER:N	1:E:63:GLN:O	2.32	0.62
1:F:89:LEU:C	1:F:90:GLU:HG2	2.20	0.62
1:B:79:PRO:O	1:B:80:ASN:CB	2.47	0.62
1:B:19:GLY:O	1:B:35:LYS:HE2	2.00	0.62
1:E:112:LYS:CE	2:E:141:SCR:O53	2.47	0.62
1:C:26:LEU:HD23	1:C:30:THR:HG21	1.80	0.62
1:B:37:ARG:O	1:B:42:ILE:HD11	1.99	0.62
1:F:114:ASN:OD1	1:F:116:ARG:HB2	1.99	0.62
1:E:24:ARG:HD3	1:E:26:LEU:CD1	2.29	0.62
1:H:85:PHE:HA	1:H:98:ILE:O	2.00	0.62
1:D:47:ALA:HB3	1:D:55:TYR:HB2	1.82	0.62
1:H:24:ARG:HD3	1:H:26:LEU:CD1	2.30	0.62
1:B:112:LYS:HA	2:B:141:SCR:O42	1.98	0.62
1:D:118:LYS:NZ	2:D:141:SCR:O44	2.31	0.62
1:A:84:LEU:CB	1:A:100:LYS:HE2	2.29	0.62
1:H:108:PHE:HB2	1:H:123:THR:CB	2.30	0.61
1:A:84:LEU:HB2	1:A:100:LYS:CD	2.29	0.61
1:G:110:GLY:O	1:G:111:LEU:HD23	2.01	0.61
1:C:68:ASP:OD1	1:C:72:LEU:HB3	2.00	0.61
1:E:89:LEU:HG	1:E:90:GLU:N	2.15	0.61
1:H:108:PHE:HB2	1:H:123:THR:HB	1.83	0.61
1:H:112:LYS:HE2	2:H:141:SCR:H61	1.82	0.61
1:C:118:LYS:HE2	2:C:141:SCR:O92	2.00	0.61
1:E:125:PHE:CD1	1:E:126:GLY:N	2.69	0.61
1:D:16:CYS:HB2	1:D:132:PHE:CE1	2.35	0.61
1:A:84:LEU:HB2	1:A:100:LYS:HE2	1.81	0.61
1:B:86:LEU:HD21	1:B:100:LYS:HG2	1.81	0.61
1:E:28:ASP:OD1	1:E:28:ASP:N	2.34	0.61
1:H:54:VAL:HG12	1:H:55:TYR:H	1.64	0.61
1:C:102:HIS:O	1:C:105:LYS:HB2	2.00	0.61
1:E:86:LEU:HD21	1:E:100:LYS:HD2	1.83	0.61
1:D:112:LYS:CE	2:D:141:SCR:O92	2.48	0.61
1:A:131:LEU:HD13	1:A:131:LEU:N	2.16	0.61
1:F:112:LYS:HE2	2:F:141:SCR:O92	2.00	0.60
1:H:53:GLU:HG2	1:H:84:LEU:CD2	2.26	0.60
1:E:86:LEU:HD21	1:E:100:LYS:CE	2.30	0.60
1:C:13:LEU:HD21	1:C:37:ARG:HG3	1.83	0.60
1:D:78:THR:HG23	1:D:79:PRO:HD2	1.81	0.60
1:H:89:LEU:C	1:H:90:GLU:HG2	2.21	0.60
1:D:112:LYS:HA	2:D:141:SCR:O42	2.02	0.60
1:E:39:ASP:O	1:E:42:ILE:HG12	2.01	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:131:LEU:HD22	1:C:131:LEU:N	2.10	0.60
1:E:112:LYS:NZ	2:E:141:SCR:O53	2.33	0.60
1:A:24:ARG:CD	1:A:26:LEU:CD1	2.79	0.60
1:D:64:PHE:CZ	1:D:79:PRO:HG3	2.36	0.60
1:E:39:ASP:OD1	1:E:39:ASP:C	2.38	0.60
1:A:114:ASN:ND2	1:A:116:ARG:HH12	1.96	0.60
2:C:141:SCR:S2	2:C:141:SCR:O84	2.60	0.60
1:B:80:ASN:O	1:B:83:CYS:N	2.33	0.60
1:G:122:ARG:HD2	2:G:141:SCR:O94	2.01	0.60
1:F:112:LYS:HG3	1:F:118:LYS:HG3	1.83	0.60
1:E:87:GLU:CG	1:E:88:ARG:N	2.64	0.59
1:E:94:TYR:C	1:E:95:ASN:HD22	2.05	0.59
1:G:104:GLU:HB2	1:G:105:LYS:HE2	1.83	0.59
1:D:123:THR:HA	1:D:127:GLN:OE1	2.03	0.59
1:A:99:SER:OG	1:A:102:HIS:HD2	1.84	0.59
1:G:89:LEU:O	1:G:90:GLU:HG2	2.03	0.59
1:C:107:TRP:CZ2	1:C:121:PRO:HG3	2.37	0.59
1:C:55:TYR:CE2	1:C:84:LEU:HD11	2.37	0.59
1:A:99:SER:OG	1:A:102:HIS:CD2	2.55	0.59
1:F:49:GLU:HB2	1:F:84:LEU:HD21	1.84	0.59
1:B:89:LEU:C	1:B:89:LEU:HD12	2.22	0.59
1:C:111:LEU:HD23	1:C:117:SER:HA	1.83	0.59
1:D:53:GLU:OE1	1:D:100:LYS:HD2	2.01	0.59
1:A:11:PRO:C	1:A:12:LYS:HD3	2.21	0.59
1:F:125:PHE:HA	1:F:130:ILE:CD1	2.33	0.59
1:C:103:ALA:C	1:C:105:LYS:H	2.06	0.59
1:G:55:TYR:O	1:G:56:ILE:HG12	2.03	0.59
1:F:28:ASP:OD1	1:F:28:ASP:C	2.41	0.59
1:G:89:LEU:C	1:G:90:GLU:HG2	2.23	0.59
1:G:66:ALA:HB1	1:G:74:TYR:CZ	2.37	0.59
2:A:141:SCR:H162	2:A:141:SCR:O5	2.01	0.59
1:G:47:ALA:O	1:G:54:VAL:HG12	2.02	0.59
1:C:64:PHE:CE2	1:C:79:PRO:HD3	2.38	0.59
1:C:26:LEU:HD23	1:C:30:THR:CG2	2.33	0.58
1:E:26:LEU:N	1:E:26:LEU:HD22	2.18	0.58
1:F:49:GLU:CG	1:F:84:LEU:HD21	2.31	0.58
2:D:141:SCR:H15	2:D:141:SCR:O52	2.01	0.58
1:G:18:ASN:ND2	1:G:111:LEU:O	2.36	0.58
1:D:102:HIS:O	1:D:106:HIS:N	2.36	0.58
1:H:130:ILE:HG22	1:H:131:LEU:HD13	1.85	0.58
1:B:48:ALA:HA	1:B:53:GLU:O	2.03	0.58
1:D:92:ASN:CG	1:D:92:ASN:O	2.39	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:103:ALA:C	1:C:105:LYS:N	2.57	0.58
2:B:141:SCR:C12	2:B:141:SCR:O93	2.51	0.58
1:H:55:TYR:CZ	1:H:84:LEU:CD1	2.87	0.58
1:A:108:PHE:O	1:A:123:THR:HG21	2.03	0.58
1:A:46:LEU:HD23	1:A:56:ILE:CG1	2.31	0.57
2:H:141:SCR:O81	2:H:141:SCR:O93	2.22	0.57
1:D:123:THR:HB	1:D:130:ILE:HD11	1.85	0.57
1:D:112:LYS:NZ	2:D:141:SCR:O92	2.36	0.57
1:F:98:ILE:HD11	1:F:106:HIS:HA	1.85	0.57
1:E:87:GLU:HG3	1:E:88:ARG:H	1.69	0.57
2:E:141:SCR:C11	2:E:141:SCR:S2	2.92	0.57
1:A:112:LYS:HG3	1:A:116:ARG:O	2.04	0.57
1:D:24:ARG:HG2	1:D:25:ILE:N	2.18	0.57
1:B:89:LEU:HD12	1:B:90:GLU:O	2.03	0.57
2:E:141:SCR:S4	2:E:141:SCR:H61	2.45	0.57
1:A:24:ARG:HD2	1:A:32:ASP:OD1	2.04	0.57
2:D:141:SCR:O81	2:D:141:SCR:O93	2.22	0.57
1:B:80:ASN:H	1:B:83:CYS:HB2	1.69	0.57
1:A:12:LYS:HG3	1:A:135:LEU:O	2.04	0.57
1:G:112:LYS:HE3	2:G:141:SCR:O43	2.05	0.56
1:H:98:ILE:HD11	1:H:108:PHE:CZ	2.40	0.56
1:G:48:ALA:HA	1:G:54:VAL:CG1	2.35	0.56
1:F:45:GLN:N	1:F:57:LYS:O	2.35	0.56
1:G:105:LYS:O	1:G:106:HIS:CB	2.52	0.56
1:H:124:HIS:O	1:H:125:PHE:O	2.23	0.56
2:H:141:SCR:H162	2:H:141:SCR:C1	2.34	0.56
1:D:98:ILE:HD11	1:D:106:HIS:C	2.24	0.56
1:D:48:ALA:CA	1:D:54:VAL:HG12	2.31	0.56
1:D:122:ARG:HD2	2:D:141:SCR:O81	2.06	0.56
1:C:131:LEU:C	1:C:131:LEU:HD23	2.25	0.56
1:D:111:LEU:HD23	1:D:117:SER:HA	1.88	0.56
1:D:39:ASP:OD1	1:D:41:HIS:N	2.36	0.56
1:G:89:LEU:CD2	1:G:90:GLU:H	2.17	0.56
1:E:64:PHE:CZ	1:E:79:PRO:HG3	2.40	0.56
2:G:141:SCR:O10	2:G:141:SCR:C5	2.53	0.56
2:B:141:SCR:H161	2:B:141:SCR:O5	2.05	0.56
1:H:37:ARG:O	1:H:42:ILE:HD11	2.05	0.56
1:A:61:THR:OG1	1:A:63:GLN:HG2	2.04	0.56
2:A:141:SCR:O93	2:A:141:SCR:S11	2.63	0.56
1:B:26:LEU:HD12	1:B:27:PRO:HD2	1.88	0.56
1:C:12:LYS:CA	1:C:135:LEU:O	2.53	0.56
1:D:124:HIS:O	1:D:127:GLN:HG3	2.04	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:54:VAL:HG12	1:H:55:TYR:N	2.21	0.56
1:H:55:TYR:CZ	1:H:84:LEU:HD12	2.41	0.56
1:H:112:LYS:HE3	2:H:141:SCR:O92	2.06	0.56
1:B:87:GLU:CG	1:B:88:ARG:N	2.68	0.56
1:E:16:CYS:SG	1:E:18:ASN:HB2	2.46	0.56
2:C:141:SCR:H162	2:C:141:SCR:O73	2.04	0.56
1:F:37:ARG:NH2	1:F:135:LEU:HD13	2.20	0.56
2:D:141:SCR:S4	2:D:141:SCR:O3	2.63	0.56
1:A:108:PHE:HB2	1:A:130:ILE:HD12	1.88	0.56
1:C:104:GLU:N	1:C:104:GLU:OE2	2.38	0.55
1:C:12:LYS:CB	1:C:135:LEU:O	2.54	0.55
1:E:112:LYS:HB2	1:E:114:ASN:OD1	2.06	0.55
1:E:52:GLY:O	1:E:86:LEU:HA	2.06	0.55
1:G:102:HIS:HA	1:G:105:LYS:HE3	1.88	0.55
1:C:125:PHE:HA	1:C:130:ILE:HD12	1.87	0.55
1:C:16:CYS:HB2	1:C:132:PHE:CD2	2.42	0.55
1:H:85:PHE:CE2	1:H:109:VAL:CG2	2.89	0.55
1:H:85:PHE:CD2	1:H:109:VAL:HG21	2.42	0.55
1:B:44:LEU:HG	1:B:56:ILE:HG22	1.88	0.55
2:A:141:SCR:O62	2:A:141:SCR:O54	2.24	0.55
1:G:112:LYS:NZ	2:G:141:SCR:C14	2.69	0.55
1:G:112:LYS:NZ	2:G:141:SCR:H14	2.20	0.55
1:C:19:GLY:HA3	1:C:21:TYR:CD2	2.41	0.55
1:D:48:ALA:HA	1:D:54:VAL:CG1	2.33	0.55
1:C:19:GLY:HA3	1:C:21:TYR:CE2	2.41	0.55
1:H:13:LEU:N	1:H:13:LEU:HD13	2.21	0.55
1:D:58:SER:O	1:D:59:THR:C	2.42	0.55
2:A:141:SCR:O93	2:A:141:SCR:O84	2.25	0.55
1:B:96:THR:HG23	1:B:108:PHE:HE2	1.72	0.55
1:D:80:ASN:HB2	1:D:83:CYS:SG	2.47	0.55
1:G:102:HIS:O	1:G:105:LYS:HG2	2.06	0.55
1:G:52:GLY:O	1:G:86:LEU:HA	2.06	0.55
1:A:26:LEU:CD1	1:A:26:LEU:H	2.20	0.55
1:C:60:GLU:N	1:C:60:GLU:CD	2.59	0.55
1:H:89:LEU:HD12	1:H:93:HIS:HA	1.89	0.55
1:H:89:LEU:CD1	1:H:93:HIS:HA	2.37	0.55
1:E:123:THR:HB	1:E:130:ILE:HD11	1.89	0.55
1:C:48:ALA:O	1:C:50:SER:N	2.40	0.55
1:C:124:HIS:HD1	1:C:124:HIS:C	2.10	0.55
2:E:141:SCR:O44	2:E:141:SCR:H61	2.06	0.55
1:A:89:LEU:O	1:A:90:GLU:OE1	2.25	0.54
1:B:22:PHE:O	1:B:23:LEU:C	2.42	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:H:71:GLY:HA3	1:H:121:PRO:CD	2.37	0.54
1:E:122:ARG:HD2	2:E:141:SCR:O92	2.06	0.54
1:H:112:LYS:CE	2:H:141:SCR:C6	2.82	0.54
1:C:104:GLU:N	1:C:104:GLU:CD	2.58	0.54
1:A:124:HIS:O	1:A:127:GLN:HG3	2.08	0.54
1:A:105:LYS:O	1:A:106:HIS:CB	2.51	0.54
2:G:141:SCR:C16	2:G:141:SCR:O63	2.56	0.54
1:D:37:ARG:O	1:D:37:ARG:CG	2.54	0.54
1:E:12:LYS:HA	1:E:135:LEU:O	2.08	0.54
1:B:80:ASN:C	1:B:82:GLU:N	2.59	0.54
1:E:24:ARG:HD2	1:E:32:ASP:OD1	2.08	0.54
1:B:15:TYR:HB2	1:B:22:PHE:HE1	1.72	0.54
1:H:25:ILE:O	1:H:25:ILE:HG22	2.07	0.54
1:E:55:TYR:CZ	1:E:84:LEU:HD22	2.43	0.54
1:A:123:THR:HB	1:A:130:ILE:HD11	1.90	0.54
1:F:49:GLU:CB	1:F:84:LEU:HD21	2.38	0.53
1:G:59:THR:HG22	1:G:60:GLU:N	2.22	0.53
1:C:78:THR:O	1:C:78:THR:HG22	2.08	0.53
2:G:141:SCR:H5	2:G:141:SCR:O10	2.08	0.53
1:D:24:ARG:HH21	1:D:34:THR:HB	1.72	0.53
1:E:55:TYR:CE1	1:E:84:LEU:HD13	2.44	0.53
1:E:27:PRO:C	1:E:29:GLY:N	2.60	0.53
1:E:49:GLU:N	1:E:53:GLU:O	2.41	0.53
1:A:102:HIS:CA	1:A:104:GLU:OE1	2.54	0.53
1:A:86:LEU:HD21	1:A:100:LYS:HB2	1.89	0.53
1:G:36:ASP:O	1:G:38:SER:N	2.42	0.53
1:G:68:ASP:OD1	1:G:72:LEU:HB3	2.08	0.53
1:F:18:ASN:HD22	1:F:129:ALA:HA	1.73	0.53
1:H:123:THR:HB	1:H:130:ILE:HD11	1.91	0.53
1:A:89:LEU:HD21	1:A:93:HIS:HA	1.90	0.53
1:G:61:THR:OG1	1:G:63:GLN:HG3	2.09	0.53
1:C:45:GLN:C	1:C:46:LEU:HD23	2.30	0.52
1:E:43:GLN:HE22	1:E:60:GLU:CG	2.22	0.52
1:A:49:GLU:HG3	1:A:55:TYR:CE2	2.44	0.52
1:G:66:ALA:CB	1:G:74:TYR:CE1	2.92	0.52
1:A:131:LEU:CD1	1:A:131:LEU:N	2.72	0.52
1:B:68:ASP:HB2	1:B:72:LEU:HB3	1.90	0.52
1:D:105:LYS:O	1:D:106:HIS:HB2	2.09	0.52
1:A:14:LEU:HD22	1:A:44:LEU:HD23	1.91	0.52
1:B:87:GLU:HG3	1:B:88:ARG:H	1.73	0.52
2:B:141:SCR:H1	2:B:141:SCR:O51	2.08	0.52
1:G:53:GLU:HB3	1:G:86:LEU:HD23	1.91	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:114:ASN:OD1	1:F:116:ARG:CB	2.58	0.52
1:H:71:GLY:HA3	1:H:121:PRO:HD2	1.90	0.52
1:H:68:ASP:HB2	1:H:72:LEU:H	1.73	0.52
1:H:24:ARG:CD	1:H:26:LEU:CD1	2.87	0.52
2:A:141:SCR:O62	2:A:141:SCR:O52	2.28	0.52
2:D:141:SCR:H5	2:D:141:SCR:O91	2.09	0.52
1:E:104:GLU:H	1:E:104:GLU:CD	2.13	0.52
1:H:85:PHE:CD2	1:H:109:VAL:CG2	2.93	0.52
1:B:15:TYR:HB2	1:B:22:PHE:CE1	2.44	0.52
1:A:78:THR:HG23	1:A:79:PRO:O	2.09	0.52
1:F:112:LYS:CE	2:F:141:SCR:O92	2.57	0.52
1:C:31:VAL:HG11	1:C:73:LEU:HD23	1.91	0.52
1:G:80:ASN:ND2	1:G:83:CYS:SG	2.83	0.52
1:H:80:ASN:HD22	1:H:80:ASN:C	2.13	0.52
1:G:21:TYR:CE1	1:G:35:LYS:HD3	2.44	0.52
1:A:112:LYS:HB3	2:A:141:SCR:O42	2.10	0.51
1:G:24:ARG:HD2	1:G:26:LEU:HD22	1.92	0.51
2:H:141:SCR:O2	2:H:141:SCR:H111	2.10	0.51
1:H:37:ARG:C	1:H:42:ILE:HD11	2.30	0.51
2:E:141:SCR:O44	2:E:141:SCR:C6	2.59	0.51
2:H:141:SCR:O53	2:H:141:SCR:H62	2.09	0.51
1:B:118:LYS:NZ	1:B:127:GLN:HE22	2.07	0.51
1:H:55:TYR:CE1	1:H:84:LEU:HD12	2.46	0.51
1:C:131:LEU:CD2	1:C:131:LEU:O	2.58	0.51
1:G:131:LEU:HD13	1:G:131:LEU:N	2.24	0.51
2:D:141:SCR:O84	2:D:141:SCR:H3	2.11	0.51
1:F:56:ILE:O	1:F:56:ILE:HG22	2.10	0.51
1:C:45:GLN:O	1:C:56:ILE:HA	2.10	0.51
1:H:44:LEU:HG	1:H:56:ILE:CG2	2.41	0.51
1:D:124:HIS:C	1:D:124:HIS:ND1	2.64	0.51
1:B:49:GLU:O	1:B:49:GLU:HG2	2.10	0.51
1:B:102:HIS:CB	1:B:105:LYS:HG3	2.38	0.51
1:D:15:TYR:HB2	1:D:135:LEU:HD11	1.91	0.51
1:H:95:ASN:HB2	1:H:132:PHE:O	2.11	0.51
1:H:12:LYS:HB2	1:H:135:LEU:O	2.11	0.51
1:H:84:LEU:O	1:H:100:LYS:HB3	2.10	0.51
1:A:22:PHE:O	1:A:33:GLY:HA2	2.11	0.51
1:G:17:SER:HB3	1:G:94:TYR:CE2	2.46	0.51
1:G:112:LYS:HE2	2:G:141:SCR:C15	2.40	0.51
1:C:71:GLY:O	1:C:120:GLY:N	2.41	0.51
1:F:28:ASP:OD1	1:F:30:THR:OG1	2.28	0.51
1:A:39:ASP:OD1	1:A:42:ILE:HG23	2.11	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:84:LEU:O	1:C:100:LYS:HB2	2.10	0.51
2:A:141:SCR:O54	2:A:141:SCR:O63	2.29	0.50
1:H:124:HIS:C	1:H:125:PHE:O	2.48	0.50
1:G:57:LYS:CG	1:G:57:LYS:O	2.58	0.50
1:D:92:ASN:C	1:D:93:HIS:CD2	2.85	0.50
1:E:27:PRO:C	1:E:29:GLY:H	2.13	0.50
1:F:92:ASN:O	1:F:93:HIS:CB	2.42	0.50
1:D:39:ASP:HB3	1:D:42:ILE:HG23	1.92	0.50
1:E:71:GLY:HA3	1:E:120:GLY:HA3	1.92	0.50
1:D:45:GLN:C	1:D:46:LEU:HD23	2.32	0.50
1:A:24:ARG:HB3	1:A:32:ASP:OD1	2.11	0.50
1:E:68:ASP:HB2	1:E:72:LEU:H	1.75	0.50
1:H:39:ASP:HB3	1:H:42:ILE:HG12	1.94	0.50
1:E:87:GLU:CG	1:E:88:ARG:H	2.24	0.50
1:G:37:ARG:HG2	1:G:37:ARG:O	2.11	0.50
2:E:141:SCR:H112	2:E:141:SCR:C2	2.41	0.50
1:C:44:LEU:HG	1:C:56:ILE:HG22	1.94	0.50
1:H:105:LYS:O	1:H:106:HIS:CB	2.59	0.50
1:E:68:ASP:HB3	1:E:70:ASP:N	2.25	0.50
1:A:112:LYS:HE3	2:A:141:SCR:O52	2.11	0.50
1:F:13:LEU:O	1:F:134:PRO:HA	2.11	0.50
1:C:131:LEU:CD2	1:C:131:LEU:C	2.79	0.50
1:H:49:GLU:HB2	1:H:53:GLU:O	2.12	0.50
1:C:25:ILE:HG23	1:C:31:VAL:HG22	1.93	0.50
1:C:24:ARG:HB3	1:C:32:ASP:O	2.12	0.50
1:G:57:LYS:HA	1:G:64:PHE:HA	1.93	0.50
2:A:141:SCR:O52	2:A:141:SCR:H61	2.10	0.50
1:G:122:ARG:C	1:G:123:THR:O	2.47	0.50
1:A:90:GLU:HB2	1:A:94:TYR:O	2.12	0.50
2:H:141:SCR:O54	2:H:141:SCR:C5	2.60	0.50
1:G:39:ASP:O	1:G:42:ILE:CG1	2.57	0.50
1:E:24:ARG:HD3	1:E:26:LEU:HD13	1.93	0.50
1:F:49:GLU:O	1:F:50:SER:CB	2.59	0.49
2:E:141:SCR:O44	2:E:141:SCR:C5	2.58	0.49
1:H:29:GLY:O	1:H:30:THR:C	2.50	0.49
2:E:141:SCR:O24	2:E:141:SCR:C11	2.58	0.49
1:D:124:HIS:O	1:D:124:HIS:ND1	2.42	0.49
1:H:53:GLU:CG	1:H:84:LEU:HD23	2.32	0.49
1:D:24:ARG:NH2	1:D:34:THR:HB	2.27	0.49
1:E:97:TYR:N	1:E:97:TYR:CD1	2.79	0.49
1:C:28:ASP:OD1	1:C:30:THR:HB	2.13	0.49
1:F:49:GLU:HB2	1:F:84:LEU:CD2	2.43	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:19:GLY:HA3	1:D:21:TYR:HE2	1.76	0.49
1:H:112:LYS:HE3	2:H:141:SCR:C6	2.41	0.49
1:H:24:ARG:CD	1:H:26:LEU:HD13	2.41	0.49
2:G:141:SCR:O6	2:G:141:SCR:C16	2.55	0.49
1:E:48:ALA:HA	1:E:54:VAL:HG12	1.93	0.49
1:C:64:PHE:N	1:C:76:SER:O	2.38	0.49
2:A:141:SCR:O54	2:A:141:SCR:S6	2.71	0.49
1:F:24:ARG:HG3	1:F:34:THR:CG2	2.43	0.49
1:A:84:LEU:CB	1:A:100:LYS:HE3	2.43	0.49
1:D:64:PHE:CE1	1:D:78:THR:HA	2.46	0.49
1:A:123:THR:CG2	1:A:130:ILE:HD11	2.43	0.49
1:H:112:LYS:NZ	2:H:141:SCR:O52	2.45	0.49
1:A:112:LYS:HZ3	1:A:112:LYS:HB3	1.77	0.48
1:D:18:ASN:O	1:D:20:GLY:N	2.46	0.48
1:G:28:ASP:OD1	1:G:30:THR:HB	2.13	0.48
1:H:89:LEU:HD11	1:H:93:HIS:ND1	2.28	0.48
1:C:89:LEU:HD13	1:C:90:GLU:H	1.72	0.48
1:H:112:LYS:HA	2:H:141:SCR:O42	2.13	0.48
2:H:141:SCR:H1	2:H:141:SCR:H162	1.96	0.48
1:C:13:LEU:HD21	1:C:37:ARG:CG	2.43	0.48
2:E:141:SCR:O91	2:E:141:SCR:O54	2.31	0.48
1:C:89:LEU:HD13	1:C:89:LEU:C	2.32	0.48
1:G:81:GLU:CB	1:G:101:LYS:HE3	2.43	0.48
1:F:36:ASP:C	1:F:38:SER:N	2.66	0.48
1:A:120:GLY:O	1:A:122:ARG:N	2.47	0.48
1:A:71:GLY:HA3	1:A:121:PRO:HD3	1.96	0.48
1:H:14:LEU:HD12	1:H:23:LEU:HD23	1.95	0.48
1:H:24:ARG:CD	1:H:32:ASP:OD1	2.61	0.48
1:A:114:ASN:OD1	1:A:116:ARG:HB2	2.14	0.48
1:C:28:ASP:C	1:C:28:ASP:OD1	2.52	0.48
1:G:48:ALA:HA	1:G:54:VAL:HG12	1.95	0.48
1:E:11:PRO:O	1:E:12:LYS:CD	2.61	0.48
1:D:46:LEU:N	1:D:46:LEU:HD23	2.28	0.48
1:F:22:PHE:O	1:F:33:GLY:HA2	2.14	0.48
1:C:112:LYS:CB	2:C:141:SCR:O42	2.61	0.48
1:B:14:LEU:HD23	1:B:44:LEU:HD23	1.96	0.48
2:D:141:SCR:O34	2:D:141:SCR:H2	2.13	0.48
1:H:104:GLU:N	1:H:104:GLU:CD	2.67	0.48
1:D:80:ASN:HB3	1:D:82:GLU:H	1.78	0.48
1:E:85:PHE:CE2	1:E:109:VAL:CG2	2.97	0.48
1:G:124:HIS:CD2	1:G:125:PHE:N	2.82	0.47
1:C:23:LEU:HD21	1:C:65:LEU:CD1	2.43	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:124:HIS:CD2	1:F:124:HIS:O	2.67	0.47
1:G:64:PHE:CE2	1:G:79:PRO:CD	2.97	0.47
1:E:105:LYS:O	1:E:106:HIS:CB	2.57	0.47
1:B:61:THR:CB	1:B:63:GLN:HE21	2.27	0.47
1:G:120:GLY:O	1:G:123:THR:OG1	2.31	0.47
1:G:57:LYS:HD2	1:G:58:SER:C	2.34	0.47
1:G:68:ASP:OD1	1:G:72:LEU:CB	2.62	0.47
1:A:84:LEU:HB3	1:A:100:LYS:HE2	1.96	0.47
1:H:88:ARG:O	1:H:89:LEU:O	2.32	0.47
1:C:112:LYS:HG2	2:C:141:SCR:O42	2.14	0.47
1:G:48:ALA:HA	1:G:54:VAL:HG13	1.97	0.47
1:B:17:SER:OG	1:B:128:LYS:HD3	2.15	0.47
1:G:122:ARG:CD	2:G:141:SCR:O94	2.62	0.47
1:G:108:PHE:HB2	1:G:123:THR:HB	1.97	0.47
1:A:89:LEU:CD2	1:A:90:GLU:N	2.69	0.47
1:D:122:ARG:CD	2:D:141:SCR:O81	2.63	0.47
1:F:130:ILE:HD13	1:F:130:ILE:HG21	1.39	0.47
1:A:57:LYS:HG2	1:A:57:LYS:O	2.14	0.47
1:F:112:LYS:CE	2:F:141:SCR:S12	2.91	0.47
1:D:112:LYS:HE2	2:D:141:SCR:O92	2.13	0.47
1:E:44:LEU:HG	1:E:56:ILE:CG2	2.45	0.47
1:F:24:ARG:O	1:F:31:VAL:HA	2.14	0.47
1:C:31:VAL:HG11	1:C:73:LEU:CD2	2.45	0.47
1:A:96:THR:HG23	1:A:130:ILE:HG23	1.96	0.47
1:E:16:CYS:HB2	1:E:132:PHE:CE2	2.50	0.47
1:D:58:SER:O	1:D:62:GLY:N	2.45	0.47
1:A:112:LYS:HB3	1:A:112:LYS:HZ2	1.78	0.47
2:B:141:SCR:O84	2:B:141:SCR:O2	2.33	0.47
1:D:89:LEU:C	1:D:90:GLU:HG2	2.35	0.47
1:F:64:PHE:CE2	1:F:79:PRO:HB3	2.50	0.47
1:A:84:LEU:HD22	1:A:84:LEU:N	2.30	0.46
2:E:141:SCR:O44	2:E:141:SCR:H5	2.16	0.46
1:B:71:GLY:HA3	1:B:121:PRO:HD3	1.96	0.46
1:H:102:HIS:HB2	1:H:107:TRP:HB2	1.96	0.46
1:G:24:ARG:CD	1:G:26:LEU:HD22	2.45	0.46
1:D:98:ILE:CD1	1:D:103:ALA:HA	2.43	0.46
1:H:112:LYS:HB3	1:H:112:LYS:HE2	1.61	0.46
1:E:80:ASN:ND2	1:E:82:GLU:HB2	2.30	0.46
1:B:104:GLU:HG3	1:B:104:GLU:H	1.36	0.46
1:E:48:ALA:HA	1:E:53:GLU:O	2.15	0.46
1:C:124:HIS:C	1:C:124:HIS:ND1	2.67	0.46
1:H:55:TYR:HE1	1:H:83:CYS:HB2	1.80	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:124:HIS:CD2	1:F:127:GLN:HG3	2.51	0.46
1:B:119:LEU:O	1:B:120:GLY:C	2.52	0.46
1:C:122:ARG:HD3	1:C:122:ARG:HH11	1.56	0.46
1:C:17:SER:HB2	1:C:131:LEU:CD2	2.46	0.46
1:H:79:PRO:O	1:H:80:ASN:CB	2.58	0.46
1:E:85:PHE:HB3	1:E:97:TYR:HB3	1.97	0.46
1:G:116:ARG:NH1	2:G:141:SCR:O52	2.49	0.46
1:F:105:LYS:O	1:F:106:HIS:CB	2.60	0.46
1:E:125:PHE:CG	1:E:126:GLY:N	2.84	0.46
1:E:12:LYS:HA	1:E:12:LYS:HD2	1.73	0.46
1:B:56:ILE:HB	1:B:65:LEU:HB2	1.97	0.46
1:D:112:LYS:CG	2:D:141:SCR:O92	2.61	0.46
1:A:80:ASN:HD22	1:A:80:ASN:C	2.20	0.46
1:A:123:THR:CB	1:A:130:ILE:HD11	2.46	0.46
1:B:61:THR:OG1	1:B:63:GLN:NE2	2.48	0.46
1:F:118:LYS:NZ	2:F:141:SCR:O84	2.49	0.46
1:E:11:PRO:O	1:E:12:LYS:HD3	2.16	0.46
1:B:17:SER:O	1:B:128:LYS:HD2	2.16	0.46
1:C:112:LYS:HA	2:C:141:SCR:O42	2.16	0.46
1:B:99:SER:O	1:B:100:LYS:C	2.54	0.46
1:B:96:THR:HG23	1:B:108:PHE:CE2	2.51	0.46
1:E:43:GLN:O	1:E:59:THR:HB	2.15	0.46
1:G:123:THR:O	1:G:124:HIS:HB3	2.16	0.45
1:B:98:ILE:HG13	1:B:108:PHE:CE2	2.51	0.45
1:D:78:THR:HG23	1:D:79:PRO:CD	2.46	0.45
2:E:141:SCR:C13	2:E:141:SCR:O74	2.60	0.45
1:B:112:LYS:C	1:B:114:ASN:H	2.20	0.45
1:E:86:LEU:HD21	1:E:100:LYS:HD3	1.94	0.45
1:F:14:LEU:HB3	1:F:132:PHE:HD2	1.81	0.45
1:G:26:LEU:HD23	1:G:30:THR:HG21	1.90	0.45
1:H:86:LEU:N	1:H:98:ILE:O	2.50	0.45
1:F:124:HIS:HD2	1:F:124:HIS:O	1.98	0.45
1:F:94:TYR:CE1	1:F:133:LEU:HB2	2.51	0.45
1:D:36:ASP:O	1:D:38:SER:N	2.49	0.45
1:F:112:LYS:HE3	2:F:141:SCR:O74	2.16	0.45
1:A:128:LYS:HA	1:A:131:LEU:HD22	1.97	0.45
1:H:118:LYS:HE2	1:H:122:ARG:O	2.17	0.45
1:H:27:PRO:CB	1:H:61:THR:HG21	2.22	0.45
1:G:112:LYS:HZ3	2:G:141:SCR:C14	2.29	0.45
1:D:24:ARG:NH1	1:D:26:LEU:HD21	2.32	0.45
1:D:112:LYS:CE	2:D:141:SCR:S12	3.00	0.45
1:A:84:LEU:O	1:A:100:LYS:HB3	2.16	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:80:ASN:C	1:E:80:ASN:HD22	2.19	0.45
1:G:124:HIS:NE2	1:G:125:PHE:O	2.50	0.45
1:B:64:PHE:O	1:B:65:LEU:C	2.55	0.45
1:G:53:GLU:HA	1:G:85:PHE:O	2.17	0.45
1:F:110:GLY:O	1:F:118:LYS:N	2.47	0.45
1:G:89:LEU:CD2	1:G:90:GLU:N	2.77	0.45
1:H:82:GLU:HG3	1:H:101:LYS:CD	2.47	0.45
1:H:73:LEU:O	1:H:74:TYR:HB3	2.16	0.45
1:B:40:GLN:O	1:B:60:GLU:HG3	2.17	0.45
1:G:105:LYS:H	1:G:105:LYS:HG2	1.42	0.45
1:A:98:ILE:HG22	1:A:99:SER:O	2.17	0.45
1:H:80:ASN:ND2	1:H:82:GLU:CB	2.80	0.45
1:E:66:ALA:HB2	1:E:83:CYS:SG	2.57	0.45
1:A:123:THR:HG22	1:A:130:ILE:HD11	1.99	0.45
1:G:59:THR:CG2	1:G:60:GLU:N	2.80	0.45
1:F:17:SER:O	1:F:18:ASN:O	2.34	0.45
2:E:141:SCR:O92	2:E:141:SCR:O54	2.35	0.45
2:A:141:SCR:O74	2:A:141:SCR:O94	2.35	0.45
1:A:26:LEU:HD11	1:A:32:ASP:OD1	2.16	0.45
1:C:104:GLU:HG2	1:C:105:LYS:CD	2.47	0.45
1:B:89:LEU:CD1	1:B:90:GLU:N	2.75	0.45
1:B:68:ASP:CB	1:B:72:LEU:CB	2.95	0.45
1:H:38:SER:O	1:H:39:ASP:C	2.52	0.45
1:H:12:LYS:HD2	1:H:46:LEU:CD1	2.47	0.45
1:F:118:LYS:HD2	2:F:141:SCR:O42	2.17	0.45
1:D:127:GLN:C	1:D:129:ALA:N	2.71	0.45
1:D:122:ARG:NH2	2:D:141:SCR:O72	2.50	0.45
1:H:26:LEU:HA	1:H:26:LEU:HD12	1.72	0.45
1:A:37:ARG:HG2	1:A:37:ARG:O	2.17	0.45
1:H:40:GLN:O	1:H:60:GLU:HG3	2.16	0.45
1:G:26:LEU:O	1:G:28:ASP:N	2.50	0.44
2:B:141:SCR:H111	2:B:141:SCR:O24	2.16	0.44
1:B:68:ASP:OD2	1:B:72:LEU:HB2	2.17	0.44
1:G:18:ASN:CG	1:G:18:ASN:O	2.54	0.44
1:G:124:HIS:O	1:G:130:ILE:HD13	2.17	0.44
1:A:89:LEU:C	1:A:90:GLU:OE1	2.56	0.44
1:A:66:ALA:HB3	1:A:74:TYR:CZ	2.52	0.44
1:E:91:GLU:C	1:E:93:HIS:H	2.20	0.44
1:H:127:GLN:O	1:H:130:ILE:HB	2.17	0.44
1:B:122:ARG:HH11	2:B:141:SCR:H112	1.82	0.44
1:A:37:ARG:O	1:A:42:ILE:HD11	2.18	0.44
1:A:17:SER:O	1:A:18:ASN:C	2.55	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:44:LEU:HG	1:C:56:ILE:CG2	2.48	0.44
2:H:141:SCR:C11	2:H:141:SCR:O2	2.65	0.44
1:F:63:GLN:HB2	1:F:63:GLN:HE21	1.46	0.44
1:E:26:LEU:CB	1:E:28:ASP:OD1	2.62	0.44
1:G:102:HIS:CB	1:G:105:LYS:HG3	2.29	0.44
1:C:128:LYS:HB2	1:C:128:LYS:HZ3	1.81	0.44
1:E:130:ILE:HD13	1:E:130:ILE:N	2.32	0.44
1:H:85:PHE:CA	1:H:98:ILE:O	2.66	0.44
1:C:135:LEU:N	1:C:135:LEU:HD23	2.31	0.44
1:H:99:SER:OG	1:H:102:HIS:CD2	2.71	0.44
1:H:12:LYS:HD2	1:H:46:LEU:HD13	1.99	0.44
1:D:130:ILE:HG22	1:D:131:LEU:CD1	2.44	0.44
1:C:112:LYS:HE3	2:C:141:SCR:O62	2.17	0.44
1:A:84:LEU:HB3	1:A:100:LYS:CE	2.48	0.44
1:B:67:MET:HG2	1:B:73:LEU:CD1	2.47	0.44
1:G:112:LYS:NZ	2:G:141:SCR:O71	2.51	0.43
1:F:112:LYS:HG2	2:F:141:SCR:O92	2.18	0.43
1:H:112:LYS:CB	2:H:141:SCR:O42	2.66	0.43
1:C:28:ASP:OD1	1:C:30:THR:CB	2.65	0.43
1:A:98:ILE:HG22	1:A:99:SER:N	2.33	0.43
1:B:71:GLY:O	1:B:120:GLY:N	2.46	0.43
1:H:21:TYR:HE2	1:H:113:LYS:O	2.01	0.43
1:F:95:ASN:O	1:F:131:LEU:HD12	2.17	0.43
1:H:68:ASP:HB3	1:H:70:ASP:H	1.81	0.43
1:G:10:LYS:HA	1:G:11:PRO:HD3	1.77	0.43
1:G:124:HIS:CD2	1:G:127:GLN:HG3	2.53	0.43
1:H:58:SER:HB3	1:H:62:GLY:H	1.83	0.43
1:A:118:LYS:HE3	2:A:141:SCR:O44	2.19	0.43
2:A:141:SCR:O74	2:A:141:SCR:C13	2.66	0.43
1:F:26:LEU:HB3	1:F:27:PRO:HD2	2.00	0.43
1:A:26:LEU:HD13	1:A:26:LEU:H	1.82	0.43
1:C:104:GLU:HG2	1:C:105:LYS:HD2	1.99	0.43
1:C:64:PHE:HE2	1:C:79:PRO:HD3	1.81	0.43
2:E:141:SCR:C11	2:E:141:SCR:O2	2.67	0.43
1:B:24:ARG:NH1	1:B:32:ASP:OD2	2.52	0.43
1:G:112:LYS:CE	2:G:141:SCR:O43	2.67	0.43
1:B:89:LEU:C	1:B:89:LEU:CD1	2.87	0.43
1:G:61:THR:HG1	1:G:63:GLN:HE21	1.61	0.43
1:H:80:ASN:ND2	1:H:82:GLU:HB2	2.33	0.43
1:F:77:GLN:HG2	1:F:77:GLN:H	1.59	0.43
1:A:112:LYS:CD	1:A:116:ARG:HB3	2.49	0.43
1:C:86:LEU:HD21	1:C:100:LYS:HG3	2.01	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:130:ILE:HD12	1:C:130:ILE:HG21	1.76	0.43
1:E:100:LYS:O	1:E:103:ALA:N	2.52	0.43
1:C:39:ASP:O	1:C:42:ILE:HG12	2.19	0.43
1:C:47:ALA:O	1:C:54:VAL:HA	2.18	0.43
2:B:141:SCR:S12	2:B:141:SCR:O1	2.77	0.43
1:D:15:TYR:CB	1:D:135:LEU:HD11	2.49	0.43
1:D:94:TYR:HB3	1:D:132:PHE:O	2.19	0.43
1:C:120:GLY:N	1:C:121:PRO:HD2	2.33	0.43
1:H:95:ASN:HD22	1:H:95:ASN:HA	1.41	0.43
1:F:78:THR:C	1:F:79:PRO:O	2.53	0.43
1:C:89:LEU:C	1:C:90:GLU:HG2	2.40	0.43
1:H:108:PHE:HB2	1:H:123:THR:OG1	2.18	0.43
1:H:124:HIS:ND1	1:H:127:GLN:HG3	2.33	0.43
1:G:84:LEU:HB3	1:G:100:LYS:HD3	2.01	0.43
1:F:96:THR:OG1	1:F:130:ILE:CG2	2.66	0.43
1:H:31:VAL:HG23	1:H:74:TYR:HA	2.01	0.43
1:H:55:TYR:CZ	1:H:84:LEU:HD11	2.53	0.43
1:H:112:LYS:NZ	2:H:141:SCR:S14	2.85	0.43
1:F:50:SER:O	1:F:51:ILE:C	2.57	0.43
1:H:24:ARG:HD2	1:H:32:ASP:OD1	2.19	0.43
1:H:24:ARG:HD3	1:H:32:ASP:OD1	2.18	0.43
1:D:89:LEU:HG	1:D:90:GLU:H	1.83	0.42
1:A:24:ARG:HD3	1:A:26:LEU:CD1	2.47	0.42
1:H:130:ILE:N	1:H:130:ILE:CD1	2.82	0.42
1:E:53:GLU:CD	1:E:100:LYS:HE3	2.39	0.42
1:A:101:LYS:HD3	1:A:102:HIS:CE1	2.54	0.42
1:B:92:ASN:ND2	1:B:94:TYR:HD2	2.17	0.42
1:H:80:ASN:C	1:H:82:GLU:N	2.73	0.42
1:H:24:ARG:CZ	1:H:26:LEU:HD11	2.48	0.42
1:A:27:PRO:HA	1:A:63:GLN:HE22	1.84	0.42
1:E:122:ARG:HD3	1:E:122:ARG:HH11	1.52	0.42
2:A:141:SCR:S14	2:A:141:SCR:O62	2.77	0.42
2:F:141:SCR:S11	2:F:141:SCR:O93	2.77	0.42
1:G:80:ASN:OD1	1:G:82:GLU:HG3	2.19	0.42
1:F:49:GLU:HG2	1:F:55:TYR:CE2	2.54	0.42
1:G:21:TYR:CD1	1:G:35:LYS:HD3	2.54	0.42
1:D:96:THR:HG22	1:D:108:PHE:CD1	2.55	0.42
1:B:112:LYS:HE2	1:B:112:LYS:HB3	1.60	0.42
1:B:80:ASN:C	1:B:82:GLU:H	2.22	0.42
1:B:90:GLU:OE2	1:B:125:PHE:CZ	2.72	0.42
1:E:44:LEU:HD13	1:E:44:LEU:N	2.34	0.42
1:C:92:ASN:O	1:C:93:HIS:HB2	2.18	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:112:LYS:HE2	2:D:141:SCR:O94	2.19	0.42
1:F:123:THR:O	1:F:124:HIS:HB3	2.20	0.42
1:F:68:ASP:HB3	1:F:74:TYR:CD2	2.55	0.42
1:H:36:ASP:O	1:H:39:ASP:HB2	2.20	0.42
1:B:18:ASN:O	1:B:20:GLY:N	2.53	0.42
2:D:141:SCR:H161	2:D:141:SCR:C1	2.50	0.42
1:F:39:ASP:OD2	1:F:41:HIS:HB2	2.19	0.42
1:C:19:GLY:CA	1:C:21:TYR:CD2	3.02	0.42
1:A:44:LEU:HG	1:A:56:ILE:CG2	2.48	0.42
1:B:108:PHE:HB3	1:B:109:VAL:H	1.68	0.42
1:H:71:GLY:HA3	1:H:121:PRO:HD3	2.02	0.42
1:C:98:ILE:HG22	1:C:99:SER:O	2.19	0.42
1:D:97:TYR:N	1:D:97:TYR:CD1	2.87	0.42
1:D:18:ASN:O	1:D:19:GLY:C	2.57	0.42
2:A:141:SCR:O93	2:A:141:SCR:H3	2.19	0.42
1:E:26:LEU:O	1:E:29:GLY:N	2.45	0.42
1:G:84:LEU:HD13	1:G:84:LEU:HA	1.68	0.42
1:G:66:ALA:CB	1:G:74:TYR:CZ	3.02	0.42
1:B:98:ILE:O	1:B:99:SER:C	2.56	0.42
1:A:63:GLN:HG2	1:A:63:GLN:H	1.53	0.42
1:D:130:ILE:HD12	1:D:130:ILE:HA	1.68	0.42
1:A:12:LYS:HD2	1:A:136:PRO:HA	2.01	0.42
1:F:86:LEU:CD2	1:F:100:LYS:HD2	2.39	0.42
1:B:89:LEU:CD1	1:B:93:HIS:HA	2.47	0.42
1:G:39:ASP:OD1	1:G:41:HIS:N	2.53	0.42
1:H:80:ASN:O	1:H:81:GLU:C	2.57	0.42
1:G:93:HIS:N	1:G:93:HIS:CD2	2.87	0.42
1:G:102:HIS:HA	1:G:105:LYS:CE	2.48	0.41
1:G:38:SER:O	1:G:39:ASP:C	2.56	0.41
1:B:49:GLU:OE1	1:B:50:SER:N	2.52	0.41
1:H:21:TYR:CE2	1:H:113:LYS:O	2.73	0.41
1:G:122:ARG:HH11	1:G:122:ARG:HD3	1.68	0.41
1:A:94:TYR:CE2	1:A:133:LEU:HB2	2.55	0.41
1:B:112:LYS:C	1:B:114:ASN:N	2.74	0.41
1:E:130:ILE:HD12	1:E:130:ILE:HA	1.78	0.41
1:F:126:GLY:H	1:F:130:ILE:HD12	1.84	0.41
1:G:64:PHE:O	1:G:65:LEU:O	2.39	0.41
1:H:120:GLY:N	1:H:121:PRO:HD2	2.36	0.41
1:H:95:ASN:HB3	1:H:96:THR:H	1.59	0.41
1:C:10:LYS:N	1:C:11:PRO:HD3	2.35	0.41
1:E:67:MET:HB2	1:E:73:LEU:HD13	2.02	0.41
2:H:141:SCR:O74	2:H:141:SCR:O91	2.39	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:40:GLN:HA	1:E:43:GLN:OE1	2.20	0.41
1:H:24:ARG:CD	1:H:26:LEU:HD11	2.49	0.41
1:H:99:SER:OG	1:H:102:HIS:HD2	2.03	0.41
1:H:66:ALA:O	1:H:73:LEU:HA	2.20	0.41
1:A:112:LYS:HG2	1:A:118:LYS:HG3	2.03	0.41
1:C:54:VAL:O	1:C:84:LEU:HA	2.19	0.41
1:A:53:GLU:HG3	1:A:86:LEU:HD23	2.03	0.41
1:E:31:VAL:HG11	1:E:65:LEU:HD21	2.03	0.41
1:G:114:ASN:ND2	2:G:141:SCR:O52	2.53	0.41
1:A:78:THR:C	1:A:79:PRO:O	2.58	0.41
1:B:61:THR:HB	1:B:63:GLN:HE21	1.84	0.41
1:C:85:PHE:HB3	1:C:97:TYR:HB3	2.03	0.41
1:E:100:LYS:O	1:E:102:HIS:N	2.53	0.41
1:B:56:ILE:HD13	1:B:65:LEU:HD12	2.02	0.41
1:D:80:ASN:HB3	1:D:81:GLU:H	1.71	0.41
1:G:43:GLN:O	1:G:59:THR:HB	2.21	0.41
1:F:73:LEU:HD21	1:F:109:VAL:HG13	2.02	0.41
1:C:77:GLN:H	1:C:77:GLN:HG2	1.23	0.41
1:D:72:LEU:HD23	1:D:73:LEU:O	2.21	0.41
1:F:112:LYS:HZ2	1:F:112:LYS:HB3	1.86	0.41
1:C:108:PHE:CE2	1:C:130:ILE:HD11	2.54	0.41
1:A:46:LEU:HB3	1:A:56:ILE:CG1	2.40	0.41
1:B:19:GLY:O	1:B:35:LYS:CE	2.67	0.41
1:D:98:ILE:HG23	1:D:98:ILE:HD13	1.82	0.41
1:G:58:SER:O	1:G:62:GLY:HA2	2.20	0.41
1:E:104:GLU:OE2	1:E:105:LYS:HD2	2.21	0.41
1:B:19:GLY:C	1:B:35:LYS:HE2	2.40	0.41
1:H:44:LEU:HB3	1:H:45:GLN:H	1.75	0.41
1:E:67:MET:HB2	1:E:73:LEU:CD1	2.51	0.41
1:B:47:ALA:O	1:B:54:VAL:HA	2.21	0.41
1:G:32:ASP:OD1	1:G:32:ASP:C	2.58	0.41
1:G:124:HIS:O	1:G:130:ILE:CD1	2.69	0.41
1:B:135:LEU:HA	1:B:136:PRO:HD3	1.75	0.41
1:F:111:LEU:HD23	1:F:117:SER:CA	2.48	0.41
1:H:104:GLU:CD	1:H:104:GLU:H	2.23	0.41
1:H:37:ARG:HA	1:H:42:ILE:HD11	2.02	0.41
1:A:84:LEU:CD2	1:A:84:LEU:N	2.84	0.40
1:G:67:MET:C	1:G:68:ASP:O	2.56	0.40
1:D:127:GLN:C	1:D:129:ALA:H	2.25	0.40
1:C:98:ILE:CG2	1:C:103:ALA:HA	2.51	0.40
2:D:141:SCR:S11	2:D:141:SCR:O93	2.79	0.40
1:B:71:GLY:C	1:B:120:GLY:H	2.23	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:16:CYS:HB2	1:F:132:PHE:CE2	2.56	0.40
1:C:26:LEU:CD2	1:C:30:THR:CG2	2.97	0.40
1:F:47:ALA:O	1:F:55:TYR:CD2	2.66	0.40
1:A:31:VAL:HG23	1:A:74:TYR:HA	2.03	0.40
1:B:49:GLU:OE1	1:B:53:GLU:CB	2.68	0.40
1:H:24:ARG:HB2	1:H:42:ILE:CG2	2.51	0.40
1:G:72:LEU:HD12	1:G:119:LEU:HD21	2.03	0.40
1:D:66:ALA:O	1:D:73:LEU:HD12	2.21	0.40
1:B:26:LEU:N	1:B:30:THR:O	2.54	0.40
2:G:141:SCR:O10	2:G:141:SCR:O5	2.38	0.40
1:D:124:HIS:C	1:D:124:HIS:HD1	2.19	0.40
1:A:89:LEU:CD2	1:A:90:GLU:O	2.69	0.40
1:B:18:ASN:HB3	1:B:19:GLY:H	1.54	0.40
1:H:122:ARG:HH11	1:H:122:ARG:HD2	1.65	0.40
1:B:24:ARG:HH11	1:B:32:ASP:CG	2.24	0.40
1:D:123:THR:O	1:D:124:HIS:CB	2.40	0.40
1:D:89:LEU:O	1:D:90:GLU:HG2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	125/140 (89%)	104 (83%)	17 (14%)	4 (3%)	6	14
1	B	125/140 (89%)	112 (90%)	9 (7%)	4 (3%)	6	14
1	C	125/140 (89%)	103 (82%)	18 (14%)	4 (3%)	6	14
1	D	125/140 (89%)	104 (83%)	16 (13%)	5 (4%)	5	9
1	E	125/140 (89%)	102 (82%)	16 (13%)	7 (6%)	3	4
1	F	125/140 (89%)	105 (84%)	11 (9%)	9 (7%)	2	2
1	G	125/140 (89%)	105 (84%)	15 (12%)	5 (4%)	5	9
1	H	125/140 (89%)	92 (74%)	19 (15%)	14 (11%)	1	0
All	All	1000/1120 (89%)	827 (83%)	121 (12%)	52 (5%)	3	5

All (52) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	81	GLU
1	B	18	ASN
1	B	68	ASP
1	D	18	ASN
1	D	50	SER
1	D	88	ARG
1	D	124	HIS
1	E	18	ASN
1	E	68	ASP
1	F	18	ASN
1	F	37	ARG
1	F	50	SER
1	F	93	HIS
1	G	39	ASP
1	G	66	ALA
1	H	30	THR
1	H	45	GLN
1	H	68	ASP
1	H	89	LEU
1	H	125	PHE
1	B	81	GLU
1	C	81	GLU
1	E	41	HIS
1	E	89	LEU
1	F	51	ILE
1	F	69	THR
1	F	89	LEU
1	G	46	LEU
1	G	60	GLU
1	H	81	GLU
1	H	91	GLU
1	H	94	TYR
1	H	130	ILE
1	A	37	ARG
1	B	93	HIS
1	F	81	GLU
1	H	39	ASP
1	H	44	LEU
1	C	68	ASP
1	D	60	GLU
1	E	96	THR

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Mol	Chain	Res	Type
1	E	101	LYS
1	G	65	LEU
1	H	60	GLU
1	C	49	GLU
1	E	60	GLU
1	A	104	GLU
1	F	135	LEU
1	H	27	PRO
1	H	50	SER
1	C	51	ILE

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	101/122 (83%)	69 (68%)	32 (32%)	0	1
1	B	101/122 (83%)	68 (67%)	33 (33%)	0	1
1	C	101/122 (83%)	67 (66%)	34 (34%)	0	0
1	D	101/122 (83%)	74 (73%)	27 (27%)	1	2
1	E	101/122 (83%)	69 (68%)	32 (32%)	0	1
1	F	101/122 (83%)	71 (70%)	30 (30%)	0	1
1	G	101/122 (83%)	64 (63%)	37 (37%)	0	0
1	H	101/122 (83%)	71 (70%)	30 (30%)	0	1
All	All	808/976 (83%)	553 (68%)	255 (32%)	0	1

All (255) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LYS
1	A	13	LEU
1	A	14	LEU
1	A	17	SER
1	A	23	LEU
1	A	26	LEU
1	A	34	THR

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Mol	Chain	Res	Type
1	A	39	ASP
1	A	44	LEU
1	A	46	LEU
1	A	49	GLU
1	A	53	GLU
1	A	57	LYS
1	A	60	GLU
1	A	65	LEU
1	A	72	LEU
1	A	73	LEU
1	A	77	GLN
1	A	78	THR
1	A	80	ASN
1	A	84	LEU
1	A	86	LEU
1	A	90	GLU
1	A	98	ILE
1	A	100	LYS
1	A	104	GLU
1	A	108	PHE
1	A	112	LYS
1	A	116	ARG
1	A	118	LYS
1	A	130	ILE
1	A	131	LEU
1	B	12	LYS
1	B	14	LEU
1	B	26	LEU
1	B	30	THR
1	B	37	ARG
1	B	44	LEU
1	B	49	GLU
1	B	53	GLU
1	B	57	LYS
1	B	59	THR
1	B	61	THR
1	B	65	LEU
1	B	67	MET
1	B	69	THR
1	B	70	ASP
1	B	72	LEU
1	B	73	LEU

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Mol	Chain	Res	Type
1	B	76	SER
1	B	78	THR
1	B	80	ASN
1	B	84	LEU
1	B	89	LEU
1	B	96	THR
1	B	100	LYS
1	B	101	LYS
1	B	104	GLU
1	B	105	LYS
1	B	122	ARG
1	B	127	GLN
1	B	128	LYS
1	B	130	ILE
1	B	131	LEU
1	B	135	LEU
1	C	12	LYS
1	C	13	LEU
1	C	14	LEU
1	C	16	CYS
1	C	17	SER
1	C	26	LEU
1	C	34	THR
1	C	43	GLN
1	C	44	LEU
1	C	46	LEU
1	C	53	GLU
1	C	57	LYS
1	C	65	LEU
1	C	67	MET
1	C	69	THR
1	C	72	LEU
1	C	77	GLN
1	C	78	THR
1	C	80	ASN
1	C	86	LEU
1	C	87	GLU
1	C	89	LEU
1	C	92	ASN
1	C	96	THR
1	C	100	LYS
1	C	101	LYS

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Mol	Chain	Res	Type
1	C	104	GLU
1	C	105	LYS
1	C	116	ARG
1	C	122	ARG
1	C	128	LYS
1	C	130	ILE
1	C	131	LEU
1	C	135	LEU
1	D	13	LEU
1	D	14	LEU
1	D	26	LEU
1	D	35	LYS
1	D	42	ILE
1	D	44	LEU
1	D	46	LEU
1	D	49	GLU
1	D	59	THR
1	D	60	GLU
1	D	63	GLN
1	D	65	LEU
1	D	68	ASP
1	D	73	LEU
1	D	77	GLN
1	D	78	THR
1	D	84	LEU
1	D	86	LEU
1	D	98	ILE
1	D	105	LYS
1	D	112	LYS
1	D	116	ARG
1	D	118	LYS
1	D	122	ARG
1	D	124	HIS
1	D	130	ILE
1	D	131	LEU
1	E	14	LEU
1	E	17	SER
1	E	30	THR
1	E	37	ARG
1	E	43	GLN
1	E	44	LEU
1	E	46	LEU

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Mol	Chain	Res	Type
1	E	57	LYS
1	E	59	THR
1	E	63	GLN
1	E	65	LEU
1	E	67	MET
1	E	68	ASP
1	E	69	THR
1	E	73	LEU
1	E	76	SER
1	E	78	THR
1	E	80	ASN
1	E	84	LEU
1	E	86	LEU
1	E	95	ASN
1	E	96	THR
1	E	97	TYR
1	E	100	LYS
1	E	101	LYS
1	E	104	GLU
1	E	116	ARG
1	E	118	LYS
1	E	122	ARG
1	E	130	ILE
1	E	131	LEU
1	E	135	LEU
1	F	13	LEU
1	F	14	LEU
1	F	24	ARG
1	F	26	LEU
1	F	30	THR
1	F	34	THR
1	F	39	ASP
1	F	44	LEU
1	F	49	GLU
1	F	54	VAL
1	F	65	LEU
1	F	68	ASP
1	F	69	THR
1	F	76	SER
1	F	77	GLN
1	F	78	THR
1	F	82	GLU

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Mol	Chain	Res	Type
1	F	84	LEU
1	F	86	LEU
1	F	89	LEU
1	F	92	ASN
1	F	94	TYR
1	F	96	THR
1	F	98	ILE
1	F	105	LYS
1	F	116	ARG
1	F	118	LYS
1	F	122	ARG
1	F	128	LYS
1	F	131	LEU
1	G	13	LEU
1	G	14	LEU
1	G	25	ILE
1	G	26	LEU
1	G	28	ASP
1	G	30	THR
1	G	35	LYS
1	G	38	SER
1	G	43	GLN
1	G	44	LEU
1	G	46	LEU
1	G	49	GLU
1	G	53	GLU
1	G	54	VAL
1	G	57	LYS
1	G	59	THR
1	G	61	THR
1	G	63	GLN
1	G	65	LEU
1	G	67	MET
1	G	68	ASP
1	G	72	LEU
1	G	73	LEU
1	G	78	THR
1	G	80	ASN
1	G	82	GLU
1	G	84	LEU
1	G	89	LEU
1	G	96	THR

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Mol	Chain	Res	Type
1	G	105	LYS
1	G	112	LYS
1	G	116	ARG
1	G	118	LYS
1	G	122	ARG
1	G	124	HIS
1	G	128	LYS
1	G	131	LEU
1	H	12	LYS
1	H	13	LEU
1	H	24	ARG
1	H	26	LEU
1	H	30	THR
1	H	34	THR
1	H	43	GLN
1	H	44	LEU
1	H	53	GLU
1	H	54	VAL
1	H	55	TYR
1	H	58	SER
1	H	65	LEU
1	H	67	MET
1	H	70	ASP
1	H	73	LEU
1	H	77	GLN
1	H	78	THR
1	H	80	ASN
1	H	84	LEU
1	H	86	LEU
1	H	89	LEU
1	H	95	ASN
1	H	96	THR
1	H	100	LYS
1	H	104	GLU
1	H	125	PHE
1	H	128	LYS
1	H	131	LEU
1	H	136	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	GLN
1	A	80	ASN
1	A	102	HIS
1	A	124	HIS
1	A	127	GLN
1	B	63	GLN
1	B	80	ASN
1	B	93	HIS
1	B	124	HIS
1	B	127	GLN
1	C	63	GLN
1	C	77	GLN
1	C	93	HIS
1	C	102	HIS
1	D	63	GLN
1	D	93	HIS
1	D	106	HIS
1	E	43	GLN
1	E	63	GLN
1	E	80	ASN
1	E	95	ASN
1	E	102	HIS
1	F	18	ASN
1	F	63	GLN
1	F	124	HIS
1	G	63	GLN
1	G	80	ASN
1	G	93	HIS
1	G	95	ASN
1	G	106	HIS
1	G	124	HIS
1	G	127	GLN
1	H	43	GLN
1	H	63	GLN
1	H	80	ASN
1	H	95	ASN
1	H	102	HIS

5.3.3 RNA ⓘ

There are no RNA chains 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	SCR	A	141	-	56,56,56	3.19	11 (19%)	92,92,92	1.75	16 (17%)
2	SCR	B	141	-	56,56,56	2.89	11 (19%)	92,92,92	1.47	12 (13%)
2	SCR	C	141	-	56,56,56	2.89	10 (17%)	92,92,92	1.75	13 (14%)
2	SCR	D	141	-	56,56,56	3.07	9 (16%)	92,92,92	1.56	14 (15%)
2	SCR	E	141	-	56,56,56	2.90	10 (17%)	92,92,92	1.69	14 (15%)
2	SCR	F	141	-	56,56,56	2.73	10 (17%)	92,92,92	1.46	13 (14%)
2	SCR	G	141	-	56,56,56	2.53	8 (14%)	92,92,92	1.90	19 (20%)
2	SCR	H	141	-	56,56,56	3.20	10 (17%)	92,92,92	1.68	15 (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	SCR	A	141	-	-	0/49/88/88	0/2/2/2
2	SCR	B	141	-	-	0/49/88/88	0/2/2/2
2	SCR	C	141	-	-	0/49/88/88	0/2/2/2
2	SCR	D	141	-	-	0/49/88/88	0/2/2/2
2	SCR	E	141	-	-	0/49/88/88	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SCR	F	141	-	-	0/49/88/88	0/2/2/2
2	SCR	G	141	-	-	0/49/88/88	0/2/2/2
2	SCR	H	141	-	-	0/49/88/88	0/2/2/2

All (79) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	141	SCR	O91-S12	-13.33	1.39	1.60
2	B	141	SCR	O91-S12	-11.26	1.42	1.60
2	H	141	SCR	O4-S4	-11.18	1.42	1.60
2	H	141	SCR	O2-S2	-10.90	1.43	1.60
2	H	141	SCR	O91-S12	-10.77	1.43	1.60
2	D	141	SCR	O2-S2	-10.60	1.43	1.60
2	A	141	SCR	O4-S4	-10.56	1.43	1.60
2	F	141	SCR	O91-S12	-10.38	1.44	1.60
2	E	141	SCR	O4-S4	-10.32	1.44	1.60
2	C	141	SCR	O91-S12	-10.31	1.44	1.60
2	E	141	SCR	O91-S12	-10.05	1.44	1.60
2	D	141	SCR	O91-S12	-10.03	1.44	1.60
2	D	141	SCR	O4-S4	-9.70	1.45	1.60
2	F	141	SCR	O4-S4	-9.61	1.45	1.60
2	D	141	SCR	O3-S3	-9.60	1.45	1.60
2	A	141	SCR	O71-S13	-9.54	1.45	1.60
2	C	141	SCR	O4-S4	-9.44	1.45	1.60
2	C	141	SCR	O71-S13	-9.29	1.45	1.60
2	H	141	SCR	O71-S13	-9.21	1.45	1.60
2	E	141	SCR	O71-S13	-8.92	1.46	1.60
2	B	141	SCR	O71-S13	-8.91	1.46	1.60
2	B	141	SCR	O4-S4	-8.80	1.46	1.60
2	G	141	SCR	O4-S4	-8.64	1.46	1.60
2	G	141	SCR	O2-S2	-8.17	1.47	1.60
2	E	141	SCR	O3-S3	-8.14	1.47	1.60
2	C	141	SCR	O2-S2	-8.09	1.47	1.60
2	B	141	SCR	O2-S2	-8.04	1.47	1.60
2	H	141	SCR	O3-S3	-7.92	1.47	1.60
2	F	141	SCR	O71-S13	-7.74	1.48	1.60
2	G	141	SCR	O71-S13	-7.63	1.48	1.60
2	D	141	SCR	O71-S13	-7.53	1.48	1.60
2	G	141	SCR	O3-S3	-7.33	1.48	1.60
2	C	141	SCR	O3-S3	-7.29	1.48	1.60
2	G	141	SCR	O91-S12	-7.25	1.48	1.60
2	B	141	SCR	O3-S3	-7.23	1.48	1.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	141	SCR	O3-S3	-7.13	1.49	1.60
2	F	141	SCR	O2-S2	-7.08	1.49	1.60
2	A	141	SCR	O2-S2	-7.02	1.49	1.60
2	F	141	SCR	O3-S3	-6.78	1.49	1.60
2	E	141	SCR	O2-S2	-6.30	1.50	1.60
2	D	141	SCR	O6-S6	-4.04	1.43	1.57
2	A	141	SCR	O91-C13	-4.03	1.40	1.46
2	C	141	SCR	O51-S14	-3.71	1.44	1.57
2	E	141	SCR	O81-S11	-3.66	1.44	1.57
2	G	141	SCR	O81-S11	-3.56	1.45	1.57
2	F	141	SCR	O6-S6	-3.56	1.45	1.57
2	B	141	SCR	O81-S11	-3.54	1.45	1.57
2	C	141	SCR	O81-S11	-3.53	1.45	1.57
2	H	141	SCR	O81-S11	-3.49	1.45	1.57
2	G	141	SCR	O6-S6	-3.48	1.45	1.57
2	E	141	SCR	O51-S14	-3.46	1.45	1.57
2	B	141	SCR	O51-S14	-3.36	1.45	1.57
2	H	141	SCR	O6-S6	-3.35	1.45	1.57
2	D	141	SCR	C11-C12	3.35	1.56	1.52
2	E	141	SCR	O6-S6	-3.27	1.46	1.57
2	F	141	SCR	C11-C12	3.26	1.55	1.52
2	A	141	SCR	O51-S14	-3.14	1.46	1.57
2	A	141	SCR	C11-C12	3.13	1.55	1.52
2	B	141	SCR	C11-C12	3.09	1.55	1.52
2	A	141	SCR	O81-S11	-3.09	1.46	1.57
2	B	141	SCR	O6-S6	-3.01	1.47	1.57
2	A	141	SCR	O6-S6	-2.90	1.47	1.57
2	G	141	SCR	O51-S14	-2.88	1.47	1.57
2	F	141	SCR	O81-S11	-2.83	1.47	1.57
2	C	141	SCR	O6-S6	-2.72	1.48	1.57
2	D	141	SCR	O81-S11	-2.70	1.48	1.57
2	D	141	SCR	O51-S14	-2.58	1.48	1.57
2	H	141	SCR	O51-S14	-2.55	1.48	1.57
2	F	141	SCR	O51-S14	-2.52	1.48	1.57
2	F	141	SCR	C4-C3	2.39	1.57	1.52
2	E	141	SCR	C12-C13	-2.39	1.50	1.54
2	H	141	SCR	O2-C2	-2.25	1.42	1.46
2	H	141	SCR	O4-C4	-2.23	1.42	1.46
2	A	141	SCR	O71-C14	-2.23	1.42	1.46
2	E	141	SCR	C3-C2	2.19	1.56	1.52
2	C	141	SCR	C12-C13	-2.17	1.50	1.54
2	B	141	SCR	O91-C13	-2.16	1.43	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	141	SCR	C4-C3	2.08	1.56	1.52
2	C	141	SCR	C4-C3	2.04	1.56	1.52

All (116) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	141	SCR	C4-O4-S4	7.77	131.05	118.37
2	G	141	SCR	C3-O3-S3	6.88	129.59	118.37
2	C	141	SCR	C16-O51-S14	-6.56	108.32	116.76
2	C	141	SCR	C1-O5-C5	-6.41	101.27	113.73
2	D	141	SCR	O4-C4-C3	-5.55	96.61	108.58
2	H	141	SCR	C14-O71-S13	-5.53	109.35	118.37
2	A	141	SCR	C16-O51-S14	5.46	123.78	116.76
2	E	141	SCR	C2-O2-S2	5.42	127.22	118.37
2	H	141	SCR	O2-C2-C1	-5.28	100.25	107.58
2	F	141	SCR	C16-O51-S14	5.26	123.52	116.76
2	H	141	SCR	C16-O51-S14	5.12	123.34	116.76
2	E	141	SCR	O6-C6-C5	5.11	117.53	107.81
2	A	141	SCR	C14-O71-S13	-4.99	110.23	118.37
2	D	141	SCR	C16-O51-S14	4.93	123.11	116.76
2	B	141	SCR	C2-O2-S2	4.82	126.23	118.37
2	E	141	SCR	C1-O5-C5	-4.81	104.38	113.73
2	G	141	SCR	O3-C3-C2	4.80	118.93	108.58
2	D	141	SCR	O81-C11-C12	4.67	115.62	108.30
2	G	141	SCR	O2-C2-C1	-4.61	101.18	107.58
2	E	141	SCR	O5-C5-C6	4.49	115.64	106.61
2	C	141	SCR	C12-O1-C1	4.49	129.16	117.50
2	A	141	SCR	C1-C2-C3	-4.27	102.78	110.70
2	B	141	SCR	C3-C4-C5	4.25	119.30	110.53
2	D	141	SCR	C14-O71-S13	4.23	125.28	118.37
2	G	141	SCR	O10-C15-C16	4.20	119.55	109.52
2	B	141	SCR	O81-C11-C12	4.11	114.73	108.30
2	G	141	SCR	C16-O51-S14	4.07	122.00	116.76
2	F	141	SCR	C3-C4-C5	4.06	118.89	110.53
2	H	141	SCR	O71-C14-C15	4.00	118.35	108.98
2	C	141	SCR	C4-O4-S4	3.98	124.86	118.37
2	A	141	SCR	C6-O6-S6	3.97	121.87	116.76
2	E	141	SCR	O5-C5-C4	-3.94	101.47	109.70
2	A	141	SCR	O91-C13-C14	-3.93	98.12	109.48
2	C	141	SCR	C6-C5-C4	3.73	122.34	113.30
2	E	141	SCR	C4-O4-S4	3.65	124.33	118.37
2	E	141	SCR	O5-C1-C2	3.65	116.90	109.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	141	SCR	C6-O6-S6	3.65	121.45	116.76
2	H	141	SCR	C1-O5-C5	-3.64	106.66	113.73
2	A	141	SCR	O6-C6-C5	3.61	114.67	107.81
2	A	141	SCR	O51-C16-C15	3.56	114.58	107.81
2	G	141	SCR	C11-C12-C13	-3.54	107.22	115.23
2	H	141	SCR	O4-C4-C3	-3.53	100.97	108.58
2	F	141	SCR	O71-C14-C13	3.52	116.75	108.88
2	H	141	SCR	C13-O91-S12	3.51	124.28	118.36
2	A	141	SCR	C6-C5-C4	3.50	121.78	113.30
2	G	141	SCR	C2-O2-S2	3.46	124.01	118.37
2	F	141	SCR	C3-O3-S3	3.45	124.00	118.37
2	A	141	SCR	O4-C4-C3	-3.44	101.17	108.58
2	E	141	SCR	C14-O71-S13	3.40	123.92	118.37
2	H	141	SCR	O91-C13-C14	-3.36	99.77	109.48
2	A	141	SCR	C3-O3-S3	3.35	123.83	118.37
2	A	141	SCR	O71-C14-C15	3.33	116.80	108.98
2	D	141	SCR	C1-O5-C5	-3.33	107.26	113.73
2	C	141	SCR	O6-C6-C5	3.29	114.06	107.81
2	F	141	SCR	O4-C4-C5	-3.24	101.72	108.59
2	H	141	SCR	C12-O1-C1	3.22	125.87	117.50
2	A	141	SCR	C2-O2-S2	3.21	123.61	118.37
2	D	141	SCR	C12-O1-C1	3.19	125.79	117.50
2	C	141	SCR	O10-C15-C14	3.16	108.88	103.48
2	C	141	SCR	C14-O71-S13	-3.06	113.38	118.37
2	B	141	SCR	C4-O4-S4	3.05	123.35	118.37
2	C	141	SCR	C16-C15-C14	-3.05	103.57	114.19
2	C	141	SCR	O5-C5-C6	3.01	112.67	106.61
2	A	141	SCR	O81-C11-C12	2.96	112.94	108.30
2	C	141	SCR	O71-C14-C13	2.96	115.50	108.88
2	E	141	SCR	C11-C12-C13	-2.95	108.56	115.23
2	E	141	SCR	C6-C5-C4	2.93	120.39	113.30
2	E	141	SCR	O10-C12-C13	2.89	109.66	105.08
2	A	141	SCR	C11-C12-C13	-2.89	108.69	115.23
2	B	141	SCR	C13-O91-S12	2.86	123.18	118.36
2	H	141	SCR	C11-C12-C13	-2.83	108.82	115.23
2	G	141	SCR	O81-C11-C12	-2.78	103.96	108.30
2	E	141	SCR	O4-C4-C3	2.76	114.53	108.58
2	B	141	SCR	C1-O5-C5	2.74	119.05	113.73
2	F	141	SCR	O10-C15-C14	2.73	108.15	103.48
2	H	141	SCR	O2-C2-C3	-2.72	102.72	108.58
2	B	141	SCR	O91-C13-C14	-2.70	101.68	109.48
2	F	141	SCR	C6-O6-S6	2.67	120.20	116.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	141	SCR	C12-O1-C1	2.66	124.42	117.50
2	F	141	SCR	O81-C11-C12	2.65	112.46	108.30
2	D	141	SCR	O81-S11-O83	2.64	114.71	106.29
2	A	141	SCR	O2-C2-C3	2.64	114.26	108.58
2	G	141	SCR	O4-C4-C5	2.59	114.10	108.59
2	B	141	SCR	O5-C5-C4	2.57	115.08	109.70
2	B	141	SCR	O6-C6-C5	2.57	112.69	107.81
2	D	141	SCR	C4-O4-S4	2.54	122.52	118.37
2	G	141	SCR	O91-S12-O92	2.54	116.05	105.54
2	B	141	SCR	C16-O51-S14	2.52	120.00	116.76
2	F	141	SCR	O2-C2-C1	2.47	111.00	107.58
2	F	141	SCR	C11-O81-S11	2.46	123.97	117.30
2	D	141	SCR	O2-C2-C1	-2.43	104.20	107.58
2	G	141	SCR	C3-C4-C5	-2.38	105.63	110.53
2	B	141	SCR	O5-C1-C2	2.36	114.29	109.52
2	G	141	SCR	O1-C1-C2	-2.31	103.56	107.58
2	H	141	SCR	O5-C5-C6	2.31	111.26	106.61
2	F	141	SCR	C4-O4-S4	2.29	122.11	118.37
2	H	141	SCR	O1-C1-C2	-2.26	103.64	107.58
2	G	141	SCR	O6-C6-C5	2.19	111.97	107.81
2	C	141	SCR	C2-O2-S2	2.19	121.94	118.37
2	F	141	SCR	O91-C13-C14	-2.19	103.16	109.48
2	F	141	SCR	O1-C1-C2	2.18	111.38	107.58
2	H	141	SCR	C1-C2-C3	2.17	114.73	110.70
2	G	141	SCR	O5-C1-C2	2.16	113.88	109.52
2	D	141	SCR	O10-C15-C14	2.12	107.11	103.48
2	D	141	SCR	O3-C3-C2	2.11	113.12	108.58
2	E	141	SCR	C3-O3-S3	2.10	121.80	118.37
2	A	141	SCR	C12-O1-C1	2.09	122.92	117.50
2	E	141	SCR	O6-S6-O62	2.06	112.87	106.29
2	D	141	SCR	O83-S11-O82	-2.06	104.45	112.89
2	G	141	SCR	O71-C14-C15	-2.06	104.15	108.98
2	G	141	SCR	O1-C12-O10	2.05	117.00	110.59
2	G	141	SCR	C1-C2-C3	-2.05	106.90	110.70
2	D	141	SCR	O1-C1-C2	-2.05	104.01	107.58
2	D	141	SCR	O44-S4-O4	2.05	115.83	107.33
2	C	141	SCR	O2-C2-C1	-2.04	104.74	107.58
2	H	141	SCR	O4-C4-C5	-2.04	104.26	108.59

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.