



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:29 PM GMT

PDB ID : 1FUG
Title : S-ADENOSYLMETHIONINE SYNTHETASE
Authors : Fu, Z.; Markham, G.D.; Takusagawa, F.
Deposited on : 1996-02-25
Resolution : 3.20 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

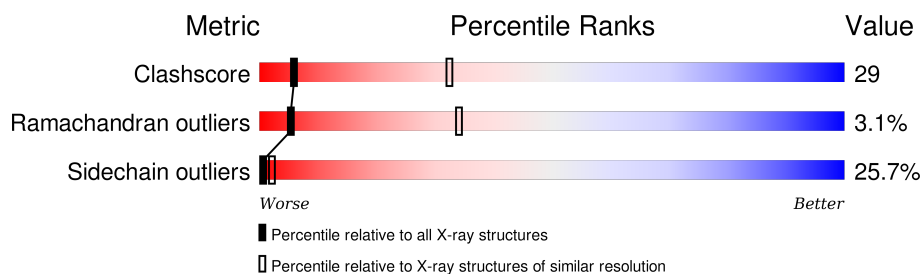
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	383	
1	B	383	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 7170 atoms, of which 1286 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 S-ADENOSYLMETHIONINE SYNTHETASE.

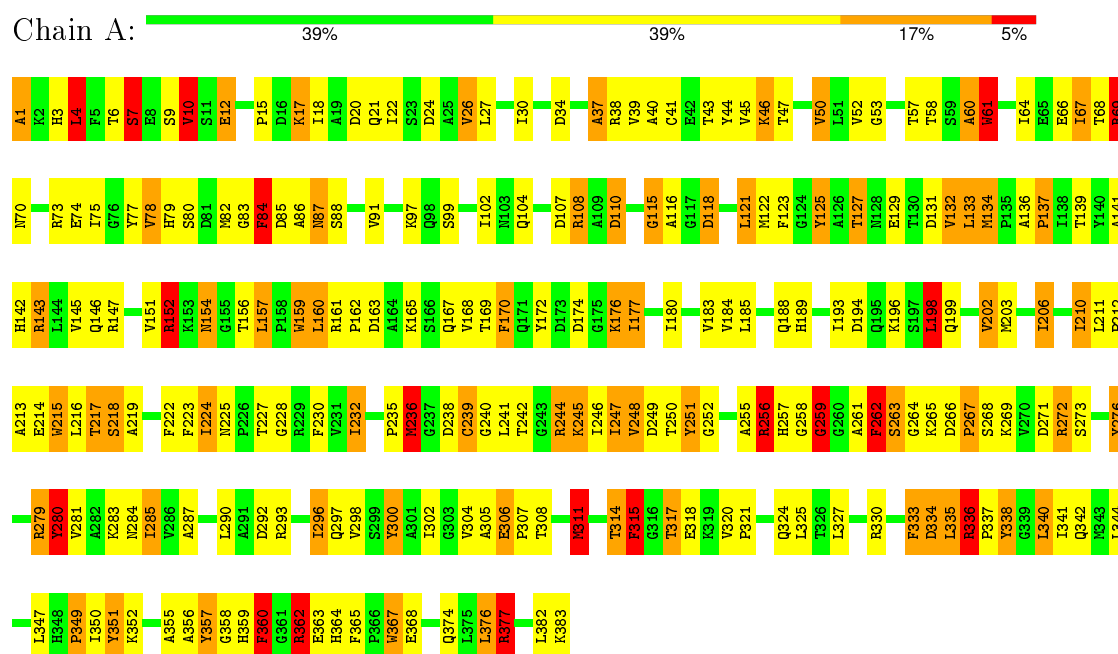
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	383	Total	C	H	N	O	S	0	0	0
			3585	1856	643	503	570	13			
1	B	383	Total	C	H	N	O	S	0	0	0
			3585	1856	643	503	570	13			

3 Residue-property plots

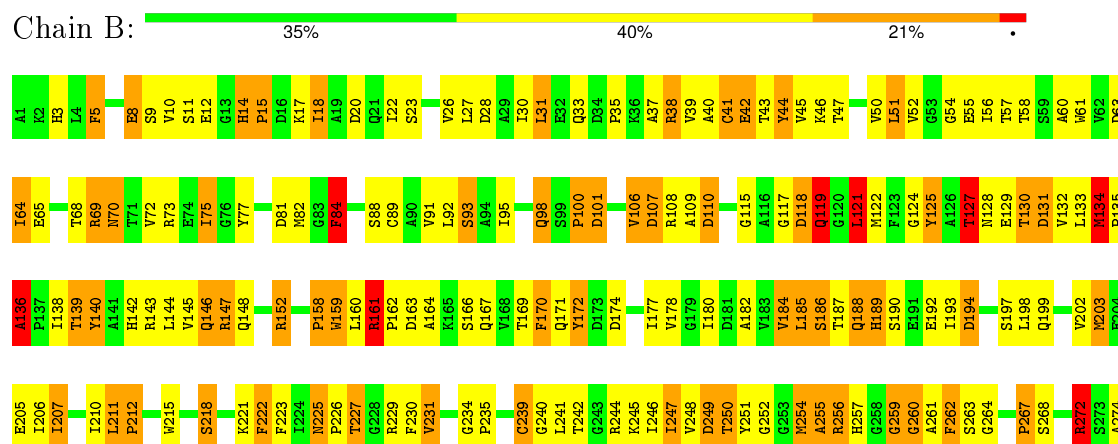
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: S-ADENOSYLMETHIONINE SYNTHETASE



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K342	K343	K344	K347	K348	K349	K350	K351	K352	K353	K354	K355	K356	K357	K358	K359	K360	K361	K362	K363	K364	K367	K368	K369	K372	K373	K374	L375	L376	K377	K378	K379	K380	K381	L382	K383	K375	K376	K377	K378	K379	K380	K381	L382	K383	K375	K376	K377	K378	K379	K380	K381	K382	K383	K384	K385	K386	K387	K388	K389	K390	K391	K392	K393	K394	K395	K396	K397	K398	K399	K400	K401	K402	K403	K404	K405	K406	K407	K408	K409	K410	K411	K412	K413	K414	K415	K416	K417	K418	K419	K420	K421	K422	K423	K424	K425	K426	K427	K428	K429	K430	K431	K432	K433	K434	K435	K436	K437	K438	K439	K440	K441	K442	K443	K444	K445	K446	K447	K448	K449	K450	K451	K452	K453	K454	K455	K456	K457	K458	K459	K460	K461	K462	K463	K464	K465	K466	K467	K468	K469	K470	K471	K472	K473	K474	K475	K476	K477	K478	K479	K480	K481	K482	K483	K484	K485	K486	K487	K488	K489	K490	K491	K492	K493	K494	K495	K496	K497	K498	K499	K500	K501	K502	K503	K504	K505	K506	K507	K508	K509	K510	K511	K512	K513	K514	K515	K516	K517	K518	K519	K520	K521	K522	K523	K524	K525	K526	K527	K528	K529	K530	K531	K532	K533	K534	K535	K536	K537	K538	K539	K540	K541	K542	K543	K544	K545	K546	K547	K548	K549	K550	K551	K552	K553	K554	K555	K556	K557	K558	K559	K560	K561	K562	K563	K564	K565	K566	K567	K568	K569	K570	K571	K572	K573	K574	K575	K576	K577	K578	K579	K580	K581	K582	K583	K584	K585	K586	K587	K588	K589	K590	K591	K592	K593	K594	K595	K596	K597	K598	K599	K600	K601	K602	K603	K604	K605	K606	K607	K608	K609	K610	K611	K612	K613	K614	K615	K616	K617	K618	K619	K620	K621	K622	K623	K624	K625	K626	K627	K628	K629	K630	K631	K632	K633	K634	K635	K636	K637	K638	K639	K640	K641	K642	K643	K644	K645	K646	K647	K648	K649	K650	K651	K652	K653	K654	K655	K656	K657	K658	K659	K660	K661	K662	K663	K664	K665	K666	K667	K668	K669	K670	K671	K672	K673	K674	K675	K676	K677	K678	K679	K680	K681	K682	K683	K684	K685	K686	K687	K688	K689	K690	K691	K692	K693	K694	K695	K696	K697	K698	K699	K700	K701	K702	K703	K704	K705	K706	K707	K708	K709	K710	K711	K712	K713	K714	K715	K716	K717	K718	K719	K720	K721	K722	K723	K724	K725	K726	K727	K728	K729	K730	K731	K732	K733	K734	K735	K736	K737	K738	K739	K740	K741	K742	K743	K744	K745	K746	K747	K748	K749	K750	K751	K752	K753	K754	K755	K756	K757	K758	K759	K760	K761	K762	K763	K764	K765	K766	K767	K768	K769	K770	K771	K772	K773	K774	K775	K776	K777	K778	K779	K780	K781	K782	K783	K784	K785	K786	K787	K788	K789	K790	K791	K792	K793	K794	K795	K796	K797	K798	K799	K800	K801	K802	K803	K804	K805	K806	K807	K808	K809	K810	K811	K812	K813	K814	K815	K816	K817	K818	K819	K820	K821	K822	K823	K824	K825	K826	K827	K828	K829	K830	K831	K832	K833	K834	K835	K836	K837	K838	K839	K840	K841	K842	K843	K844	K845	K846	K847	K848	K849	K850	K851	K852	K853	K854	K855	K856	K857	K858	K859	K860	K861	K862	K863	K864	K865	K866	K867	K868	K869	K870	K871	K872	K873	K874	K875	K876	K877	K878	K879	K880	K881	K882	K883	K884	K885	K886	K887	K888	K889	K890	K891	K892	K893	K894	K895	K896	K897	K898	K899	K900	K901	K902	K903	K904	K905	K906	K907	K908	K909	K910	K911	K912	K913	K914	K915	K916	K917	K918	K919	K920	K921	K922	K923	K924	K925	K926	K927	K928	K929	K930	K931	K932	K933	K934	K935	K936	K937	K938	K939	K940	K941	K942	K943	K944	K945	K946	K947	K948	K949	K950	K951	K952	K953	K954	K955	K956	K957	K958	K959	K960	K961	K962	K963	K964	K965	K966	K967	K968	K969	K970	K971	K972	K973	K974	K975	K976	K977	K978	K979	K980	K981	K982	K983	K984	K985	K986	K987	K988	K989	K990	K991	K992	K993	K994	K995	K996	K997	K998	K999	K1000
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4 Data and refinement statistics

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

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	121.00Å 121.00Å 171.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 3.20	Depositor
% Data completeness (in resolution range)	92.4 (8.00-3.20)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.210 , 0.290	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7170	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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	0.95	0/3001	1.93	76/4068 (1.9%)
1	B	0.92	1/3001 (0.0%)	1.97	79/4068 (1.9%)
All	All	0.94	1/6002 (0.0%)	1.95	155/8136 (1.9%)

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	19
1	B	0	18
All	All	0	37

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	260	GLY	CA-C	5.17	1.60	1.51

All (155) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	279	ARG	NE-CZ-NH1	13.88	127.24	120.30
1	B	279	ARG	NE-CZ-NH2	-12.27	114.17	120.30
1	A	244	ARG	NE-CZ-NH2	-11.39	114.61	120.30
1	A	244	ARG	NE-CZ-NH1	11.14	125.87	120.30
1	B	272	ARG	NE-CZ-NH2	-10.51	115.05	120.30
1	B	377	ARG	NE-CZ-NH1	10.25	125.42	120.30
1	A	279	ARG	NE-CZ-NH1	10.07	125.33	120.30
1	B	330	ARG	NE-CZ-NH1	9.67	125.14	120.30
1	B	82	MET	CA-CB-CG	9.66	129.72	113.30
1	B	152	ARG	NE-CZ-NH2	-9.53	115.53	120.30
1	B	351	TYR	CB-CG-CD1	-9.30	115.42	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	161	ARG	NE-CZ-NH1	8.89	124.75	120.30
1	A	61	TRP	CD1-CG-CD2	8.88	113.41	106.30
1	B	367	TRP	CD1-CG-CD2	8.67	113.24	106.30
1	A	239	CYS	CA-CB-SG	-8.62	98.48	114.00
1	B	259	GLY	CA-C-N	8.62	133.44	116.20
1	A	215	TRP	CD1-CG-CD2	8.46	113.07	106.30
1	B	326	THR	CA-CB-CG2	8.46	124.24	112.40
1	A	61	TRP	CB-CG-CD1	-8.42	116.06	127.00
1	B	250	THR	CA-C-N	8.34	135.56	117.20
1	A	280	TYR	CB-CG-CD2	-8.25	116.05	121.00
1	A	236	MET	CG-SD-CE	-8.23	87.02	100.20
1	A	351	TYR	CB-CG-CD1	-8.07	116.16	121.00
1	A	61	TRP	CE2-CD2-CG	-8.03	100.87	107.30
1	B	101	ASP	CA-C-N	8.02	134.85	117.20
1	A	61	TRP	CG-CD2-CE3	7.98	141.08	133.90
1	B	261	ALA	O-C-N	7.91	135.36	122.70
1	B	159	TRP	CB-CG-CD1	-7.86	116.79	127.00
1	B	261	ALA	CA-C-N	-7.81	100.02	117.20
1	A	168	VAL	CG1-CB-CG2	-7.79	98.43	110.90
1	A	336	ARG	NE-CZ-NH2	-7.74	116.43	120.30
1	B	159	TRP	CE2-CD2-CG	-7.70	101.14	107.30
1	A	134	MET	CG-SD-CE	-7.60	88.03	100.20
1	A	255	ALA	CB-CA-C	-7.48	98.88	110.10
1	B	367	TRP	CE2-CD2-CG	-7.39	101.39	107.30
1	B	159	TRP	CG-CD2-CE3	7.38	140.54	133.90
1	A	7	SER	CA-CB-OG	-7.36	91.32	111.20
1	B	159	TRP	CD1-CG-CD2	7.25	112.10	106.30
1	B	375	LEU	CA-CB-CG	7.25	131.98	115.30
1	B	215	TRP	CD1-CG-CD2	7.15	112.02	106.30
1	A	26	VAL	CG1-CB-CG2	-7.11	99.53	110.90
1	A	10	VAL	CA-CB-CG2	-7.09	100.27	110.90
1	B	147	ARG	NE-CZ-NH1	7.08	123.84	120.30
1	A	382	LEU	CA-CB-CG	6.95	131.28	115.30
1	B	215	TRP	CE2-CD2-CG	-6.88	101.80	107.30
1	B	108	ARG	NE-CZ-NH1	6.88	123.74	120.30
1	B	100	PRO	N-CA-C	6.83	129.85	112.10
1	A	215	TRP	CE2-CD2-CG	-6.81	101.86	107.30
1	B	338	TYR	CB-CG-CD1	-6.74	116.96	121.00
1	A	152	ARG	CA-CB-CG	6.74	128.22	113.40
1	A	159	TRP	CD1-CG-CD2	6.74	111.69	106.30
1	B	84	PHE	CB-CG-CD1	6.73	125.51	120.80
1	A	152	ARG	NE-CZ-NH2	-6.68	116.96	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	119	GLN	CA-CB-CG	6.64	128.02	113.40
1	B	61	TRP	CE2-CD2-CG	-6.59	102.03	107.30
1	B	61	TRP	CD1-CG-CD2	6.47	111.48	106.30
1	B	249	ASP	CA-CB-CG	6.38	127.44	113.40
1	A	293	ARG	NE-CZ-NH2	-6.33	117.14	120.30
1	A	60	ALA	CB-CA-C	-6.30	100.65	110.10
1	B	260	GLY	N-CA-C	6.30	128.84	113.10
1	A	159	TRP	CE2-CD2-CG	-6.28	102.28	107.30
1	B	69	ARG	NE-CZ-NH1	6.26	123.43	120.30
1	A	227	THR	CA-C-N	-6.26	103.68	116.20
1	B	364	HIS	N-CA-CB	6.25	121.86	110.60
1	A	67	ILE	CG1-CB-CG2	-6.24	97.68	111.40
1	A	37	ALA	CB-CA-C	-6.20	100.80	110.10
1	B	314	THR	N-CA-C	-6.20	94.27	111.00
1	A	259	GLY	CA-C-N	6.13	128.46	116.20
1	B	256	ARG	NE-CZ-NH1	6.12	123.36	120.30
1	A	215	TRP	CG-CD1-NE1	-6.12	103.98	110.10
1	A	272	ARG	NE-CZ-NH1	6.12	123.36	120.30
1	A	78	VAL	O-C-N	-6.07	112.99	122.70
1	A	198	LEU	CA-CB-CG	6.02	129.14	115.30
1	A	125	TYR	CB-CG-CD1	-6.01	117.39	121.00
1	A	271	ASP	CB-CA-C	-5.99	98.43	110.40
1	B	152	ARG	NE-CZ-NH1	5.98	123.29	120.30
1	A	146	GLN	CA-CB-CG	-5.95	100.32	113.40
1	B	326	THR	CA-CB-OG1	-5.93	96.54	109.00
1	A	61	TRP	CG-CD1-NE1	-5.92	104.18	110.10
1	B	367	TRP	CG-CD2-CE3	5.92	139.23	133.90
1	B	167	GLN	CB-CA-C	-5.92	98.57	110.40
1	B	367	TRP	CG-CD1-NE1	-5.91	104.19	110.10
1	B	250	THR	O-C-N	-5.90	113.26	122.70
1	B	259	GLY	C-N-CA	-5.89	109.92	122.30
1	A	262	PHE	CB-CG-CD2	-5.89	116.68	120.80
1	B	167	GLN	N-CA-CB	5.81	121.06	110.60
1	A	367	TRP	CE2-CD2-CG	-5.81	102.66	107.30
1	B	383	LYS	N-CA-C	5.80	126.65	111.00
1	A	215	TRP	CB-CG-CD1	-5.78	119.49	127.00
1	B	8	GLU	CA-CB-CG	5.77	126.09	113.40
1	A	269	LYS	N-CA-C	-5.76	95.45	111.00
1	A	217	THR	CA-CB-CG2	5.74	120.44	112.40
1	B	212	PRO	CA-C-N	-5.70	104.66	117.20
1	B	136	ALA	N-CA-C	5.64	126.24	111.00
1	B	51	LEU	CA-CB-CG	5.64	128.27	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	147	ARG	NE-CZ-NH2	-5.63	117.48	120.30
1	A	152	ARG	NE-CZ-NH1	5.63	123.12	120.30
1	A	314	THR	N-CA-C	-5.63	95.80	111.00
1	B	121	LEU	CA-CB-CG	5.62	128.22	115.30
1	A	334	ASP	CA-C-N	-5.60	104.87	117.20
1	A	290	LEU	CA-CB-CG	5.57	128.11	115.30
1	B	128	ASN	N-CA-CB	5.56	120.61	110.60
1	B	134	MET	CG-SD-CE	-5.56	91.30	100.20
1	A	306	GLU	CB-CA-C	5.54	121.49	110.40
1	A	338	TYR	CB-CG-CD1	-5.53	117.68	121.00
1	B	186	SER	N-CA-CB	5.52	118.78	110.50
1	B	239	CYS	CA-CB-SG	-5.52	104.07	114.00
1	A	1	ALA	N-CA-C	-5.51	96.13	111.00
1	A	133	LEU	N-CA-CB	-5.50	99.41	110.40
1	A	223	PHE	CB-CG-CD2	-5.49	116.96	120.80
1	A	377	ARG	NE-CZ-NH1	5.49	123.04	120.30
1	B	101	ASP	O-C-N	-5.47	113.95	122.70
1	B	61	TRP	N-CA-C	-5.47	96.24	111.00
1	A	143	ARG	NE-CZ-NH1	5.45	123.03	120.30
1	B	351	TYR	CA-C-N	5.45	129.19	117.20
1	A	69	ARG	NE-CZ-NH2	-5.44	117.58	120.30
1	B	360	PHE	N-CA-C	-5.43	96.34	111.00
1	A	202	VAL	CG1-CB-CG2	-5.42	102.23	110.90
1	A	367	TRP	CD1-CG-CD2	5.41	110.63	106.30
1	A	236	MET	CA-CB-CG	5.41	122.49	113.30
1	A	263	SER	CB-CA-C	-5.39	99.86	110.10
1	A	357	TYR	CA-CB-CG	-5.39	103.16	113.40
1	B	125	TYR	CB-CG-CD2	-5.38	117.78	121.00
1	B	44	TYR	N-CA-C	-5.36	96.52	111.00
1	B	255	ALA	CB-CA-C	-5.35	102.08	110.10
1	A	315	PHE	CA-C-N	-5.34	105.52	116.20
1	B	259	GLY	CA-C-O	-5.34	110.99	120.60
1	B	84	PHE	CA-CB-CG	5.33	126.70	113.90
1	A	50	VAL	N-CA-C	-5.33	96.60	111.00
1	A	176	LYS	CA-C-N	-5.32	105.49	117.20
1	B	146	GLN	N-CA-C	-5.30	96.69	111.00
1	B	14	HIS	CA-CB-CG	5.23	122.49	113.60
1	A	215	TRP	CG-CD2-CE3	5.22	138.60	133.90
1	A	362	ARG	N-CA-C	-5.21	96.92	111.00
1	B	20	ASP	N-CA-CB	-5.20	101.24	110.60
1	A	133	LEU	CB-CA-C	5.19	120.06	110.20
1	A	4	LEU	CA-CB-CG	5.17	127.20	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	127	THR	CA-C-N	5.15	128.53	117.20
1	B	267	PRO	N-CD-CG	-5.15	95.47	103.20
1	A	311	MET	CA-CB-CG	5.15	122.05	113.30
1	A	256	ARG	NE-CZ-NH1	5.15	122.87	120.30
1	B	377	ARG	NE-CZ-NH2	-5.13	117.73	120.30
1	A	185	LEU	CB-CG-CD1	-5.12	102.29	111.00
1	B	336	ARG	CB-CG-CD	5.12	124.92	111.60
1	B	260	GLY	CA-C-O	5.11	129.79	120.60
1	A	232	ILE	N-CA-C	-5.10	97.24	111.00
1	B	250	THR	N-CA-C	5.10	124.77	111.00
1	A	108	ARG	NE-CZ-NH1	5.09	122.85	120.30
1	A	279	ARG	NE-CZ-NH2	-5.05	117.77	120.30
1	B	42	GLU	CB-CA-C	-5.05	100.30	110.40
1	B	311	MET	CA-CB-CG	5.05	121.89	113.30
1	B	88	SER	N-CA-CB	-5.05	102.93	110.50
1	A	102	ILE	O-C-N	-5.05	114.63	122.70
1	B	229	ARG	NE-CZ-NH2	-5.01	117.79	120.30
1	A	102	ILE	CA-C-N	5.01	128.22	117.20

There are no chirality outliers.

All (37) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	110	ASP	Peptide
1	A	125	TYR	Sidechain
1	A	132	VAL	Mainchain
1	A	143	ARG	Sidechain
1	A	152	ARG	Mainchain
1	A	170	PHE	Sidechain
1	A	251	TYR	Sidechain
1	A	256	ARG	Sidechain
1	A	276	TYR	Sidechain
1	A	280	TYR	Sidechain
1	A	300	TYR	Sidechain
1	A	315	PHE	Sidechain
1	A	351	TYR	Sidechain
1	A	357	TYR	Sidechain
1	A	360	PHE	Sidechain
1	A	362	ARG	Sidechain
1	A	44	TYR	Sidechain
1	A	69	ARG	Sidechain
1	A	84	PHE	Sidechain

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Mol	Chain	Res	Type	Group
1	B	106	VAL	Peptide
1	B	107	ASP	Peptide
1	B	140	TYR	Sidechain
1	B	170	PHE	Peptide
1	B	172	TYR	Sidechain
1	B	260	GLY	Peptide
1	B	272	ARG	Sidechain
1	B	276	TYR	Sidechain
1	B	280	TYR	Sidechain
1	B	294	CYS	Peptide
1	B	300	TYR	Sidechain
1	B	315	PHE	Sidechain
1	B	336	ARG	Sidechain
1	B	338	TYR	Sidechain
1	B	359	HIS	Sidechain
1	B	363	GLU	Mainchain
1	B	38	ARG	Sidechain
1	B	84	PHE	Sidechain

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	2942	643	2908	168	0
1	B	2942	643	2908	182	0
All	All	5884	1286	5816	336	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 29.

All (336) 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:45:VAL:HG12	1:B:50:VAL:HG13	1.54	0.87
1:A:30:ILE:HG12	1:A:58:THR:HG21	1.59	0.85
1:B:251:TYR:HB2	1:B:255:ALA:HB2	1.60	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203:MET:SD	1:B:222:PHE:HE2	2.01	0.83
1:B:134:MET:SD	1:B:279:ARG:NH2	2.51	0.83
1:B:276:TYR:HE1	1:B:359:HIS:HB3	1.45	0.81
1:B:45:VAL:HG22	1:B:240:GLY:HA3	1.64	0.80
1:B:152:ARG:HB3	1:B:160:LEU:HD11	1.64	0.80
1:A:121:LEU:HB2	1:A:297:GLN:HE22	1.47	0.79
1:A:134:MET:SD	1:A:279:ARG:NH2	2.55	0.79
1:A:15:PRO:HB2	1:A:240:GLY:HA3	1.64	0.78
1:A:177:ILE:HG13	1:A:215:TRP:CZ2	2.21	0.76
1:B:65:GLU:HG3	1:B:91:VAL:HG11	1.67	0.76
1:B:278:ALA:HA	1:B:281:VAL:HG23	1.69	0.75
1:B:134:MET:HG3	1:B:135:PRO:HD2	1.66	0.75
1:B:138:ILE:HB	1:B:252:GLY:HA2	1.69	0.74
1:A:292:ASP:HB2	1:A:317:THR:HG22	1.67	0.74
1:B:203:MET:SD	1:B:222:PHE:CE2	2.81	0.72
1:B:130:THR:OG1	1:B:136:ALA:HB3	1.93	0.69
1:A:21:GLN:HE21	1:A:352:LYS:HE3	1.56	0.68
1:B:72:VAL:HG11	1:B:84:PHE:CE1	2.28	0.68
1:A:281:VAL:HG21	1:A:296:ILE:HD12	1.76	0.68
1:A:266:ASP:HB2	1:A:267:PRO:HD2	1.74	0.68
1:A:157:LEU:HD23	1:A:159:TRP:CZ2	2.30	0.67
1:B:11:SER:HA	1:B:357:TYR:HE1	1.60	0.66
1:B:134:MET:HA	1:B:279:ARG:HH22	1.60	0.66
1:B:5:PHE:CE2	1:B:254:MET:SD	2.88	0.66
1:A:161:ARG:O	1:A:188:GLN:HB3	1.96	0.66
1:B:122:MET:HG2	1:B:274:ALA:HB3	1.78	0.66
1:A:157:LEU:HD22	1:A:160:LEU:HD21	1.78	0.66
1:A:203:MET:SD	1:A:222:PHE:CD2	2.89	0.65
1:A:46:LYS:HB3	1:A:239:CYS:SG	2.37	0.65
1:B:138:ILE:HB	1:B:252:GLY:CA	2.27	0.65
1:A:9:SER:HB2	1:A:141:ALA:HB1	1.78	0.65
1:A:10:VAL:HG23	1:A:165:LYS:HA	1.79	0.65
1:B:129:GLU:HG2	1:B:135:PRO:HA	1.78	0.65
1:B:147:ARG:HD3	1:B:206:ILE:HA	1.78	0.64
1:A:261:ALA:HB3	1:A:265:LYS:NZ	2.13	0.64
1:B:50:VAL:HG23	1:B:89:CYS:SG	2.37	0.64
1:A:297:GLN:HB3	1:A:311:MET:HB3	1.80	0.64
1:A:30:ILE:HG21	1:A:39:VAL:HG11	1.79	0.64
1:B:145:VAL:HG12	1:B:206:ILE:HD11	1.80	0.63
1:A:262:PHE:O	1:A:272:ARG:HD3	1.99	0.63
1:A:245:LYS:HA	1:A:245:LYS:HE3	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:351:TYR:O	1:B:354:THR:HG22	1.99	0.62
1:B:11:SER:HB3	1:B:164:ALA:HB3	1.81	0.62
1:B:134:MET:SD	1:B:279:ARG:CZ	2.87	0.62
1:B:109:ALA:HB2	1:B:115:GLY:HA2	1.82	0.62
1:B:42:GLU:HA	1:B:242:THR:HG21	1.82	0.62
1:A:83:GLY:HA3	1:A:236:MET:HB2	1.80	0.62
1:B:72:VAL:HG11	1:B:84:PHE:HE1	1.65	0.62
1:A:37:ALA:CB	1:A:58:THR:HG22	2.30	0.62
1:B:357:TYR:H	1:B:362:ARG:NH2	1.97	0.61
1:A:172:TYR:HE1	1:A:177:ILE:HG22	1.65	0.61
1:A:12:GLU:HB2	1:A:21:GLN:OE1	2.01	0.61
1:B:298:VAL:HG13	1:B:310:ILE:HG22	1.83	0.61
1:A:52:VAL:HG11	1:A:64:ILE:HG21	1.83	0.61
1:A:314:THR:HB	1:A:317:THR:OG1	2.01	0.60
1:B:15:PRO:HA	1:B:18:ILE:HG22	1.82	0.60
1:A:167:GLN:NE2	1:B:119:GLN:HG2	2.17	0.60
1:A:30:ILE:CG2	1:A:39:VAL:HG11	2.32	0.60
1:A:37:ALA:HB2	1:A:58:THR:HG22	1.83	0.60
1:A:281:VAL:CG2	1:A:296:ILE:HD12	2.32	0.59
1:B:65:GLU:HA	1:B:91:VAL:HG21	1.83	0.59
1:A:9:SER:OG	1:A:142:HIS:HD2	1.86	0.59
1:A:263:SER:HA	1:A:272:ARG:NH1	2.18	0.59
1:A:320:VAL:HB	1:A:321:PRO:HD2	1.83	0.59
1:A:247:ILE:HB	1:B:259:GLY:CA	2.33	0.59
1:B:184:VAL:HB	1:B:223:PHE:HB2	1.84	0.59
1:B:295:GLU:HG2	1:B:315:PHE:HE1	1.68	0.59
1:A:46:LYS:HB3	1:A:239:CYS:HG	1.67	0.59
1:B:319:LYS:N	1:B:319:LYS:HD2	2.18	0.59
1:B:185:LEU:HD21	1:B:206:ILE:HD13	1.84	0.58
1:A:272:ARG:HG2	1:A:276:TYR:CE2	2.38	0.58
1:A:256:ARG:HH12	1:B:256:ARG:HH22	1.51	0.58
1:B:23:SER:CB	1:B:41:CYS:SG	2.92	0.58
1:A:122:MET:SD	1:A:122:MET:N	2.77	0.57
1:B:133:LEU:HD12	1:B:283:LYS:HG3	1.87	0.57
1:A:239:CYS:SG	1:A:240:GLY:N	2.76	0.57
1:A:188:GLN:HA	1:A:230:PHE:HB3	1.85	0.57
1:A:250:THR:HG22	1:A:360:PHE:CE1	2.39	0.57
1:B:10:VAL:HA	1:B:164:ALA:O	2.03	0.57
1:B:164:ALA:HA	1:B:186:SER:O	2.05	0.57
1:A:159:TRP:O	1:A:189:HIS:HA	2.04	0.57
1:A:199:GLN:HA	1:A:224:ILE:HD11	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:LYS:NZ	1:A:356:ALA:HA	2.19	0.57
1:B:23:SER:HB3	1:B:41:CYS:SG	2.45	0.56
1:A:242:THR:HG22	1:B:244:ARG:NH2	2.19	0.56
1:B:161:ARG:HE	1:B:190:SER:HA	1.70	0.56
1:B:27:LEU:HD13	1:B:31:LEU:HD21	1.87	0.56
1:A:159:TRP:CD2	1:A:193:ILE:HG13	2.40	0.56
1:B:122:MET:SD	1:B:122:MET:N	2.78	0.56
1:B:171:GLN:HG3	1:B:178:VAL:HG13	1.87	0.56
1:B:367:TRP:O	1:B:368:GLU:HG3	2.06	0.56
1:A:276:TYR:HE1	1:A:359:HIS:HD2	1.52	0.56
1:B:378:ASP:O	1:B:382:LEU:HD23	2.06	0.56
1:A:84:PHE:HD2	1:A:235:PRO:HD2	1.70	0.56
1:B:152:ARG:HH22	1:B:162:PRO:HG3	1.71	0.56
1:B:11:SER:HA	1:B:357:TYR:CE1	2.41	0.56
1:B:272:ARG:NH2	1:B:354:THR:HG21	2.21	0.56
1:A:118:ASP:HA	1:A:302:ILE:HA	1.87	0.55
1:A:37:ALA:HA	1:A:57:THR:O	2.06	0.55
1:A:285:ILE:HA	1:A:376:LEU:HD23	1.89	0.55
1:B:127:THR:HB	1:B:129:GLU:HB2	1.89	0.55
1:A:46:LYS:HA	1:A:235:PRO:HB3	1.89	0.55
1:B:335:LEU:HA	1:B:340:LEU:HD21	1.88	0.55
1:B:296:ILE:HA	1:B:311:MET:O	2.07	0.55
1:A:177:ILE:HG13	1:A:215:TRP:HZ2	1.71	0.55
1:A:121:LEU:HB2	1:A:297:GLN:NE2	2.18	0.54
1:A:276:TYR:HE1	1:A:359:HIS:CD2	2.24	0.54
1:B:18:ILE:HD11	1:B:72:VAL:HG22	1.88	0.54
1:A:308:THR:HG23	1:B:221:LYS:NZ	2.23	0.54
1:A:320:VAL:HG23	1:A:321:PRO:O	2.07	0.54
1:A:159:TRP:CZ3	1:A:198:LEU:HG	2.42	0.54
1:A:259:GLY:N	1:B:247:ILE:HG21	2.22	0.54
1:B:140:TYR:O	1:B:144:LEU:HB2	2.08	0.54
1:A:259:GLY:HA2	1:B:247:ILE:HB	1.90	0.54
1:A:156:THR:HG22	1:A:157:LEU:HD12	1.90	0.53
1:A:40:ALA:HA	1:A:264:GLY:O	2.08	0.53
1:B:56:ILE:HG13	1:B:95:ILE:HD12	1.91	0.53
1:B:272:ARG:HH21	1:B:354:THR:HG21	1.72	0.53
1:B:245:LYS:O	1:B:249:ASP:HB2	2.08	0.53
1:A:280:TYR:O	1:A:284:ASN:HB2	2.08	0.53
1:A:22:ILE:HD13	1:A:50:VAL:HG11	1.90	0.53
1:A:82:MET:O	1:A:232:ILE:HA	2.09	0.53
1:A:338:TYR:CE1	1:A:341:ILE:HD12	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:320:VAL:HG21	1:A:324:GLN:HB3	1.90	0.53
1:B:43:THR:HG22	1:B:52:VAL:HA	1.91	0.53
1:B:246:ILE:HB	1:B:257:HIS:HE1	1.74	0.53
1:A:157:LEU:HB2	1:A:160:LEU:HD11	1.91	0.53
1:B:117:GLY:O	1:B:302:ILE:HG13	2.08	0.53
1:B:278:ALA:HA	1:B:281:VAL:CG2	2.37	0.52
1:A:202:VAL:HA	1:A:206:ILE:HG12	1.90	0.52
1:B:18:ILE:HG21	1:B:235:PRO:HG3	1.91	0.52
1:A:133:LEU:HD21	1:A:287:ALA:HB2	1.92	0.52
1:B:15:PRO:O	1:B:18:ILE:HG22	2.09	0.52
1:B:41:CYS:HA	1:B:54:GLY:HA3	1.91	0.52
1:A:257:HIS:O	1:B:247:ILE:HG13	2.10	0.52
1:A:324:GLN:HG3	1:A:327:LEU:HD12	1.91	0.52
1:B:30:ILE:HG23	1:B:58:THR:HG21	1.92	0.52
1:B:57:THR:HG22	1:B:98:GLN:HB3	1.92	0.52
1:A:172:TYR:CE1	1:A:177:ILE:HG22	2.44	0.51
1:A:298:VAL:HG21	1:A:335:LEU:HD22	1.92	0.51
1:A:247:ILE:HB	1:B:259:GLY:HA2	1.92	0.51
1:A:121:LEU:HG	1:A:258:GLY:HA3	1.93	0.51
1:A:261:ALA:HB3	1:A:265:LYS:HZ1	1.76	0.51
1:A:83:GLY:CA	1:A:236:MET:HB2	2.41	0.51
1:A:300:TYR:CE1	1:A:307:PRO:HB3	2.45	0.51
1:B:44:TYR:HB3	1:B:51:LEU:HD21	1.92	0.51
1:A:3:HIS:HB2	1:A:172:TYR:HB2	1.93	0.50
1:A:242:THR:HG22	1:B:244:ARG:HH21	1.75	0.50
1:B:69:ARG:HH11	1:B:69:ARG:HG2	1.75	0.50
1:A:157:LEU:N	1:A:157:LEU:HD12	2.26	0.50
1:B:189:HIS:O	1:B:231:VAL:HG23	2.11	0.50
1:B:170:PHE:CE1	1:B:180:ILE:HD12	2.46	0.50
1:B:121:LEU:C	1:B:122:MET:SD	2.89	0.50
1:A:147:ARG:O	1:A:151:VAL:HG22	2.12	0.50
1:A:374:GLN:O	1:A:377:ARG:HG3	2.12	0.50
1:B:325:LEU:HA	1:B:328:LEU:HB3	1.93	0.50
1:B:194:ASP:O	1:B:197:SER:HB3	2.11	0.50
1:B:202:VAL:HG13	1:B:206:ILE:CG2	2.41	0.50
1:A:64:ILE:HG22	1:A:68:THR:OG1	2.12	0.50
1:A:133:LEU:HD13	1:A:283:LYS:HE3	1.94	0.50
1:B:8:GLU:HG3	1:B:248:VAL:HG11	1.94	0.50
1:A:245:LYS:HG3	1:A:248:VAL:HG11	1.95	0.49
1:A:276:TYR:HB3	1:A:367:TRP:HB3	1.95	0.49
1:A:64:ILE:HG22	1:A:68:THR:HG1	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:324:GLN:O	1:A:327:LEU:HB3	2.12	0.49
1:B:57:THR:O	1:B:58:THR:HB	2.13	0.49
1:B:161:ARG:HH22	1:B:231:VAL:HG22	1.77	0.49
1:B:144:LEU:HD11	1:B:207:ILE:HA	1.95	0.49
1:B:147:ARG:HB2	1:B:210:ILE:HD11	1.95	0.49
1:A:116:ALA:N	1:A:305:ALA:HB2	2.28	0.49
1:A:115:GLY:H	1:A:305:ALA:HB3	1.77	0.49
1:A:188:GLN:HB2	1:A:230:PHE:HD2	1.77	0.49
1:A:70:ASN:HA	1:A:73:ARG:HD2	1.94	0.49
1:A:224:ILE:HD12	1:A:225:ASN:HB2	1.94	0.48
1:B:186:SER:HA	1:B:225:ASN:HB3	1.96	0.48
1:A:70:ASN:O	1:A:73:ARG:HB2	2.13	0.48
1:A:61:TRP:CZ3	1:A:97:LYS:HD3	2.48	0.48
1:A:34:ASP:HB3	1:A:58:THR:HB	1.94	0.48
1:B:118:ASP:HA	1:B:302:ILE:HB	1.93	0.48
1:B:31:LEU:HB3	1:B:349:PRO:HD3	1.95	0.48
1:B:17:LYS:HD3	1:B:17:LYS:HA	1.77	0.48
1:A:27:LEU:HD13	1:A:39:VAL:CG2	2.43	0.48
1:B:125:TYR:HE2	1:B:135:PRO:HB3	1.79	0.48
1:B:247:ILE:HG23	1:B:252:GLY:O	2.13	0.48
1:B:124:GLY:HA3	1:B:278:ALA:O	2.14	0.48
1:B:169:THR:HB	1:B:182:ALA:HB3	1.95	0.48
1:B:17:LYS:HD2	1:B:355:ALA:O	2.13	0.48
1:B:26:VAL:HG21	1:B:64:ILE:HG22	1.96	0.48
1:B:185:LEU:HD13	1:B:207:ILE:HD12	1.94	0.47
1:B:35:PRO:HB2	1:B:347:LEU:HG	1.95	0.47
1:B:37:ALA:HB1	1:B:57:THR:O	2.14	0.47
1:A:53:GLY:HA3	1:B:44:TYR:OH	2.14	0.47
1:A:15:PRO:HG3	1:A:238:ASP:OD2	2.15	0.47
1:A:66:GLU:O	1:A:70:ASN:HB2	2.15	0.47
1:B:333:PHE:CE1	1:B:372:LYS:HD2	2.49	0.47
1:B:161:ARG:NE	1:B:190:SER:HA	2.29	0.47
1:B:262:PHE:HE1	1:B:275:ALA:HB3	1.80	0.47
1:B:138:ILE:HD11	1:B:248:VAL:O	2.15	0.47
1:A:211:LEU:HA	1:A:212:PRO:HD3	1.72	0.47
1:B:239:CYS:SG	1:B:240:GLY:N	2.87	0.47
1:A:27:LEU:HA	1:A:30:ILE:HG22	1.97	0.47
1:A:77:TYR:HB3	1:A:84:PHE:O	2.13	0.47
1:B:31:LEU:HD12	1:B:35:PRO:HA	1.96	0.47
1:B:23:SER:OG	1:B:43:THR:HG23	2.15	0.47
1:B:8:GLU:HA	1:B:166:SER:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:SER:HB2	1:A:137:PRO:HB2	1.96	0.46
1:B:277:ALA:HA	1:B:344:LEU:HD21	1.96	0.46
1:B:276:TYR:CE1	1:B:359:HIS:HB3	2.37	0.46
1:A:245:LYS:HA	1:A:245:LYS:CE	2.45	0.46
1:B:132:VAL:HG13	1:B:134:MET:H	1.80	0.46
1:A:4:LEU:O	1:B:309:SER:HB3	2.15	0.46
1:B:161:ARG:HG3	1:B:161:ARG:HH11	1.80	0.46
1:B:262:PHE:CE1	1:B:275:ALA:HB3	2.50	0.46
1:B:44:TYR:HB3	1:B:51:LEU:CD2	2.45	0.46
1:B:33:GLN:OE1	1:B:60:ALA:HA	2.15	0.46
1:B:295:GLU:HG2	1:B:315:PHE:CE1	2.48	0.46
1:B:161:ARG:NH2	1:B:231:VAL:HG22	2.31	0.46
1:B:5:PHE:CZ	1:B:254:MET:SD	3.09	0.45
1:A:9:SER:HB2	1:A:141:ALA:CB	2.46	0.45
1:B:182:ALA:HB2	1:B:221:LYS:HD3	1.98	0.45
1:B:27:LEU:O	1:B:31:LEU:HD22	2.16	0.45
1:B:277:ALA:O	1:B:280:TYR:N	2.50	0.45
1:B:31:LEU:CB	1:B:349:PRO:HD3	2.46	0.45
1:A:305:ALA:HB1	1:A:336:ARG:HB2	1.97	0.45
1:A:337:PRO:O	1:A:340:LEU:HD22	2.17	0.45
1:A:152:ARG:HD2	1:A:162:PRO:HG3	1.97	0.45
1:A:174:ASP:C	1:A:176:LYS:H	2.20	0.45
1:B:152:ARG:NH2	1:B:162:PRO:HG3	2.30	0.45
1:B:77:TYR:OH	1:B:234:GLY:HA2	2.16	0.45
1:B:357:TYR:N	1:B:362:ARG:NH2	2.63	0.45
1:B:22:ILE:HG21	1:B:68:THR:HG23	1.98	0.45
1:B:202:VAL:HA	1:B:206:ILE:HG22	1.99	0.45
1:A:320:VAL:HG22	1:A:325:LEU:HD12	1.98	0.45
1:B:127:THR:O	1:B:133:LEU:HA	2.16	0.45
1:B:131:ASP:OD1	1:B:132:VAL:HG12	2.17	0.45
1:B:328:LEU:HD13	1:B:380:ALA:HB2	1.99	0.45
1:B:91:VAL:O	1:B:92:LEU:HD23	2.17	0.45
1:A:24:ASP:OD1	1:A:264:GLY:N	2.50	0.45
1:A:338:TYR:CD1	1:A:341:ILE:HD12	2.52	0.45
1:A:261:ALA:HB3	1:A:265:LYS:HZ2	1.78	0.45
1:B:325:LEU:HB2	1:B:328:LEU:HD22	1.99	0.45
1:A:333:PHE:CD1	1:A:333:PHE:N	2.85	0.45
1:A:167:GLN:HB3	1:A:184:VAL:CG2	2.47	0.44
1:B:296:ILE:HD11	1:B:298:VAL:HG22	1.98	0.44
1:A:276:TYR:CE1	1:A:359:HIS:CD2	3.05	0.44
1:A:20:ASP:HB3	1:A:355:ALA:HB1	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:246:ILE:HD12	1:B:257:HIS:ND1	2.33	0.44
1:B:46:LYS:HB3	1:B:239:CYS:SG	2.57	0.44
1:A:306:GLU:HA	1:A:307:PRO:HD3	1.83	0.44
1:B:300:TYR:CD2	1:B:337:PRO:HG3	2.53	0.44
1:A:222:PHE:N	1:A:222:PHE:CD1	2.85	0.44
1:A:236:MET:SD	1:A:236:MET:O	2.76	0.44
1:B:280:TYR:CE2	1:B:343:MET:SD	3.11	0.44
1:B:70:ASN:O	1:B:73:ARG:HB3	2.18	0.44
1:A:74:GLU:OE2	1:A:352:LYS:HD2	2.18	0.44
1:B:225:ASN:HA	1:B:226:PRO:HD2	1.82	0.44
1:B:144:LEU:HD11	1:B:207:ILE:HG13	2.00	0.44
1:B:280:TYR:HE2	1:B:343:MET:SD	2.41	0.44
1:B:163:ASP:O	1:B:187:THR:HA	2.18	0.44
1:A:267:PRO:O	1:A:341:ILE:HD11	2.17	0.44
1:A:79:HIS:H	1:A:82:MET:HE3	1.82	0.44
1:B:247:ILE:O	1:B:250:THR:N	2.51	0.43
1:B:246:ILE:HB	1:B:257:HIS:CE1	2.53	0.43
1:B:281:VAL:HG21	1:B:296:ILE:HD12	1.99	0.43
1:B:64:ILE:HD11	1:B:93:SER:OG	2.19	0.43
1:A:210:ILE:HD13	1:A:210:ILE:HA	1.95	0.43
1:A:213:ALA:HA	1:A:216:LEU:HD12	2.00	0.43
1:B:77:TYR:CZ	1:B:162:PRO:HG2	2.51	0.43
1:B:140:TYR:HD1	1:B:143:ARG:HG3	1.82	0.43
1:B:134:MET:HG3	1:B:135:PRO:CD	2.43	0.43
1:B:140:TYR:CD1	1:B:143:ARG:HG3	2.53	0.43
1:A:127:THR:HG23	1:A:129:GLU:H	1.84	0.43
1:B:152:ARG:CB	1:B:160:LEU:HD11	2.43	0.43
1:A:69:ARG:HD2	1:A:86:ALA:O	2.19	0.43
1:B:188:GLN:HA	1:B:230:PHE:HB3	2.00	0.43
1:B:27:LEU:HD13	1:B:31:LEU:CD2	2.47	0.43
1:A:334:ASP:HB3	1:A:336:ARG:NE	2.34	0.43
1:B:251:TYR:HE1	1:B:279:ARG:HH21	1.67	0.42
1:B:139:THR:O	1:B:143:ARG:HG2	2.19	0.42
1:A:320:VAL:CG2	1:A:324:GLN:HB3	2.49	0.42
1:B:54:GLY:O	1:B:95:ILE:HA	2.19	0.42
1:A:136:ALA:N	1:A:137:PRO:HD2	2.33	0.42
1:A:64:ILE:O	1:A:64:ILE:HG22	2.19	0.42
1:B:288:ALA:HB2	1:B:373:ALA:HB1	2.01	0.42
1:B:84:PHE:CD2	1:B:235:PRO:HB2	2.54	0.42
1:A:268:SER:HB2	1:A:338:TYR:OH	2.20	0.42
1:A:64:ILE:HG21	1:A:64:ILE:HD13	1.80	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:134:MET:SD	1:B:279:ARG:NH1	2.93	0.42
1:A:266:ASP:HB2	1:A:267:PRO:CD	2.45	0.42
1:A:26:VAL:CG2	1:A:64:ILE:HG23	2.50	0.42
1:B:338:TYR:O	1:B:341:ILE:HB	2.19	0.42
1:A:308:THR:HG23	1:B:221:LYS:HZ2	1.83	0.42
1:A:61:TRP:HZ3	1:A:97:LYS:HD3	1.82	0.42
1:B:211:LEU:HA	1:B:212:PRO:HD3	1.71	0.42
1:A:115:GLY:N	1:A:305:ALA:HB3	2.34	0.42
1:A:217:THR:HB	1:A:218:SER:H	1.73	0.42
1:A:365:PHE:HB2	1:A:368:GLU:HB2	2.02	0.42
1:B:40:ALA:HA	1:B:264:GLY:O	2.20	0.42
1:B:294:CYS:SG	1:B:295:GLU:N	2.93	0.42
1:A:68:THR:HG21	1:A:91:VAL:CG2	2.50	0.42
1:A:133:LEU:N	1:A:133:LEU:HD12	2.35	0.42
1:A:183:VAL:HG11	1:A:211:LEU:HD22	2.02	0.41
1:A:18:ILE:HD11	1:A:75:ILE:HD11	2.02	0.41
1:B:46:LYS:HB3	1:B:239:CYS:HG	1.85	0.41
1:A:250:THR:HG22	1:A:360:PHE:CZ	2.55	0.41
1:B:15:PRO:O	1:B:18:ILE:N	2.54	0.41
1:A:123:PHE:CD2	1:A:297:GLN:HB2	2.55	0.41
1:A:217:THR:CB	1:A:219:ALA:H	2.34	0.41
1:A:9:SER:OG	1:A:142:HIS:HA	2.21	0.41
1:A:213:ALA:HA	1:A:216:LEU:HB2	2.03	0.41
1:A:311:MET:HG2	1:B:3:HIS:CG	2.56	0.41
1:B:10:VAL:HB	1:B:14:HIS:HB3	2.01	0.41
1:A:87:ASN:OD1	1:A:87:ASN:N	2.54	0.41
1:A:358:GLY:N	1:A:362:ARG:HH22	2.18	0.41
1:A:170:PHE:CZ	1:A:180:ILE:HD12	2.55	0.41
1:A:3:HIS:O	1:A:172:TYR:N	2.53	0.41
1:B:15:PRO:HG3	1:B:235:PRO:HA	2.03	0.41
1:B:161:ARG:NH1	1:B:161:ARG:HG3	2.36	0.41
1:B:98:GLN:HB2	1:B:98:GLN:HE21	1.80	0.41
1:A:377:ARG:HB2	1:A:383:LYS:HA	2.03	0.41
1:A:350:ILE:HD12	1:A:350:ILE:HA	1.92	0.41
1:B:12:GLU:O	1:B:75:ILE:HG21	2.20	0.41
1:A:154:ASN:N	1:A:154:ASN:HD22	2.18	0.41
1:A:6:THR:HG23	1:A:169:THR:HG22	2.02	0.41
1:B:199:GLN:O	1:B:202:VAL:HB	2.21	0.41
1:A:362:ARG:HD2	1:A:364:HIS:HE1	1.86	0.41
1:A:1:ALA:O	1:A:3:HIS:ND1	2.54	0.40
1:B:193:ILE:HD11	1:B:197:SER:OG	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:MET:SD	1:A:161:ARG:HD2	2.60	0.40
1:A:83:GLY:HA2	1:A:236:MET:HG2	2.02	0.40
1:B:288:ALA:HB3	1:B:290:LEU:HD11	2.02	0.40
1:B:279:ARG:NH1	1:B:283:LYS:NZ	2.69	0.40
1:A:304:VAL:HG12	1:A:306:GLU:O	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	381/383 (100%)	303 (80%)	69 (18%)	9 (2%)	7	43
1	B	381/383 (100%)	295 (77%)	71 (19%)	15 (4%)	4	28
All	All	762/766 (100%)	598 (78%)	140 (18%)	24 (3%)	5	34

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	47	THR
1	B	136	ALA
1	A	228	GLY
1	B	47	THR
1	B	227	THR
1	B	231	VAL
1	B	241	LEU
1	B	350	ILE
1	A	60	ALA
1	A	110	ASP
1	A	115	GLY
1	A	252	GLY
1	B	100	PRO

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Mol	Chain	Res	Type
1	B	110	ASP
1	B	218	SER
1	B	254	MET
1	B	322	SER
1	B	334	ASP
1	B	158	PRO
1	B	267	PRO
1	A	259	GLY
1	B	321	PRO
1	A	137	PRO
1	A	349	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	311/311 (100%)	236 (76%)	75 (24%)	1	3
1	B	311/311 (100%)	226 (73%)	85 (27%)	0	2
All	All	622/622 (100%)	462 (74%)	160 (26%)	0	2

All (160) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	A	7	SER
1	A	10	VAL
1	A	12	GLU
1	A	17	LYS
1	A	38	ARG
1	A	41	CYS
1	A	43	THR
1	A	45	VAL
1	A	46	LYS
1	A	61	TRP
1	A	67	ILE

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Mol	Chain	Res	Type
1	A	69	ARG
1	A	78	VAL
1	A	80	SER
1	A	84	PHE
1	A	85	ASP
1	A	87	ASN
1	A	88	SER
1	A	99	SER
1	A	104	GLN
1	A	107	ASP
1	A	108	ARG
1	A	118	ASP
1	A	121	LEU
1	A	127	THR
1	A	131	ASP
1	A	132	VAL
1	A	139	THR
1	A	145	VAL
1	A	152	ARG
1	A	154	ASN
1	A	157	LEU
1	A	160	LEU
1	A	163	ASP
1	A	177	ILE
1	A	194	ASP
1	A	196	LYS
1	A	198	LEU
1	A	206	ILE
1	A	210	ILE
1	A	214	GLU
1	A	218	SER
1	A	224	ILE
1	A	236	MET
1	A	241	LEU
1	A	244	ARG
1	A	245	LYS
1	A	246	ILE
1	A	247	ILE
1	A	248	VAL
1	A	249	ASP
1	A	251	TYR
1	A	262	PHE

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Mol	Chain	Res	Type
1	A	267	PRO
1	A	273	SER
1	A	285	ILE
1	A	296	ILE
1	A	311	MET
1	A	315	PHE
1	A	317	THR
1	A	318	GLU
1	A	330	ARG
1	A	333	PHE
1	A	335	LEU
1	A	336	ARG
1	A	340	LEU
1	A	342	GLN
1	A	344	LEU
1	A	347	LEU
1	A	349	PRO
1	A	360	PHE
1	A	363	GLU
1	A	376	LEU
1	A	377	ARG
1	B	5	PHE
1	B	9	SER
1	B	15	PRO
1	B	18	ILE
1	B	28	ASP
1	B	31	LEU
1	B	38	ARG
1	B	39	VAL
1	B	41	CYS
1	B	55	GLU
1	B	63	ASP
1	B	64	ILE
1	B	70	ASN
1	B	75	ILE
1	B	81	ASP
1	B	84	PHE
1	B	93	SER
1	B	98	GLN
1	B	101	ASP
1	B	106	VAL
1	B	107	ASP

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Mol	Chain	Res	Type
1	B	110	ASP
1	B	118	ASP
1	B	119	GLN
1	B	121	LEU
1	B	127	THR
1	B	130	THR
1	B	131	ASP
1	B	134	MET
1	B	139	THR
1	B	142	HIS
1	B	146	GLN
1	B	148	GLN
1	B	158	PRO
1	B	159	TRP
1	B	161	ARG
1	B	172	TYR
1	B	174	ASP
1	B	177	ILE
1	B	184	VAL
1	B	185	LEU
1	B	188	GLN
1	B	189	HIS
1	B	192	GLU
1	B	194	ASP
1	B	198	LEU
1	B	203	MET
1	B	205	GLU
1	B	207	ILE
1	B	211	LEU
1	B	218	SER
1	B	222	PHE
1	B	225	ASN
1	B	227	THR
1	B	247	ILE
1	B	262	PHE
1	B	263	SER
1	B	268	SER
1	B	279	ARG
1	B	281	VAL
1	B	283	LYS
1	B	284	ASN
1	B	290	LEU

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Mol	Chain	Res	Type
1	B	292	ASP
1	B	296	ILE
1	B	308	THR
1	B	310	ILE
1	B	311	MET
1	B	312	VAL
1	B	314	THR
1	B	315	PHE
1	B	317	THR
1	B	318	GLU
1	B	320	VAL
1	B	325	LEU
1	B	328	LEU
1	B	331	GLU
1	B	332	PHE
1	B	336	ARG
1	B	347	LEU
1	B	352	LYS
1	B	360	PHE
1	B	362	ARG
1	B	364	HIS
1	B	369	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	HIS
1	A	104	GLN
1	A	142	HIS
1	A	154	ASN
1	A	297	GLN
1	B	98	GLN
1	B	257	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

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

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

6.3 Carbohydrates [i](#)

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

6.4 Ligands [i](#)

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

6.5 Other polymers [i](#)

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