Analysis of NH_2-NH_3^+ molecules in CSD

Version 5.32 plus updates (Nov 2010, Feb 2011, May 2011, Aug 2011, Nov 2011)
Search criteria: 3-D coordinates determined, no powder structures, only organics

COUNT	MIN	MAX	MEAN	STD.DEV
49	1.396	1.463	1.442	0.011

NAME	DISTANCE	REFERENCE (Journal, volume, page, year)
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