SUPPLEMENTARY INFORMATION

for

Feinberg, L.F., R. Srikanth, R.W. Vachet, and J.F. Holden (submitted) "Constraints on anaerobic growth in the hyperthermophilic Archaea *Pyrobaculum islandicum* and *Pyrobaculum aerophilum*". Appl. Environ. Microbiol. (submitted 9/5/07, revised 11/5/07).

The following information is available at:

http://www.bio.umass.edu/micro/holden/suppldata-1.html

and this website is cited in the journal article.



FIG. S1. Electrospray ionization mass spectra of one of the compounds eluting at 3.2 min. The spectra identify this compound as cytidine. (a) MS/MS of m/z 244; the ion at m/z 112 is protonated cytosine, and the mass difference between m/z 244 and 112 is consistent with the neutral loss of a ribose unit. (b) Experimentally measured and theoretically predicted isotope distribution for $(M+H)^+$ of cytidine.



FIG. S2. Electrospray ionization mass spectra of one of the compounds eluting at 3.2 min. The spectra identify this compound as uridine. (a) MS/MS of m/z 243; the ion at m/z 110 is deprotonated uracil, and the mass difference between m/z 243 and 110 is consistent with the neutral loss of a ribose unit. (b) Experimentally measured and theoretically predicted isotope distribution for $(M-H)^{-}$ of uridine.



FIG. S3. Electrospray ionization mass spectra of one of the compounds eluting at 3.4 min. The spectra identify this compound as adenosine. (a) MS/MS of m/z 268; the ion at m/z 136 is protonated adenine, and the mass difference between m/z 268 and 136 is consistent with the neutral loss of a ribose unit. (b) MS³ of the product ion at m/z 136, confirming the structure of adenine (Nelson and McCloskey, 1992). (c) Experimentally measured and theoretically predicted isotope distribution for $(M+H)^+$ of adenosine.



FIG. S4. Electrospray ionization mass spectra of one of the compounds eluting at 3.4 min. The spectra identify this compound as guanosine. (a) MS/MS of m/z 284; the ion at m/z 152 is protonated guanine, and the mass difference between m/z 284 and 152 is consistent with the neutral loss of a ribose unit. (b) Experimentally measured and theoretically predicted isotope distribution for $(M+H)^+$ of guanosine.

TABLE 1. Peptide masses measured from the trypsin-digested 120-kDa protein from

thiosulfate-grown cultures compared with the predicted trypsin-digested masses of the α

subunit of thiosulfate reductase (Pisl_0266) and a solute binding protein (Pisl_0913)

Measured	Predicted	Δ m/z	Amino acid sequence			
m/z	m/z					
Thiosulfate reductase, α subunit (Pisl_0266):						
1495.6	1495.7	0.1	YSVDLEEWLWR (806-816)			
2208.8	2209.0	0.2	DYMEAAAREADVDVNDLRR (415-433)			
2413.9	2414.1	0.2	FADTYGGGVITFDNPYCTYPR (193-213)			
3013.7	3014.6	0.9	DVVVRVHVYKPVKYSVDLEEWLWR (793-816)			
3060.3	3060.4	0.1	DGTEWSQNMPVGMITEEGIFPIPNLER (711-737)			
3250.6	3250.7	0.1	IVDEAGNVRAPTVEDIYKMGGYMVLVPTGR			
			(748-777)			
3616.9	3616.9	0	VHVYKPVKYSVDLEEWLWRTIHYNSPMAR			
			(798-826)			

Predicted amino acid sequence (matches are in red): MSRFFLLMEISRREVLKAGATIGLIGGVSGILLKAAVEQSKAEAASAVTSVPSICGMCMAQCAIYIDV VDGKPVRIRPNTNAPTSAKGICARGVAGTFNAWLNPDAIKKPMARRALVDWAQGKISWEEAKRQL VTNRGKYDDMVEVDWNTAIDIIAKKLKELADNNERHAFTFLFGAWGPVASMRAGVPLMRFADTY **GGGVITFDNPYCTYPR**YLGHWLTWGHGHQAHVACIDYGEAEAVLVVRRNVIGAGVVTETWRFME **AVRRGAKLVVLSPVFDETASYADVWLPVKPGTDLAVLLAFIKYVLDNGYYMAEYLRRFTNAPFLIK** PDGLPLLASEVDWGKYGVKEPAFAYVVWDEAAGGPAPDNAAQRAALFGEYEVALKDGSVVKAKT ALQILKEWVDANLSALAOKHGVKDYMEAAAREADVDVNDLRRAAEIVAKYRAVSPIGWHDPRYS NSPQTWRAVGVLMALLGRIQQPGGLFLLTHLIMPYADVYTKVMKYTKKDVPYKTIRGLTFGEYVS ANLRGIYVIPIAPPLPGPSDRDAPPVKSLTEVWGEEAEKKGYLYPYDTVQALYESVVHGKPFKIKVV FITGSNPIPRIGNSRLVEEIFRNLELVIVHDIQFNDTTAFADVILPDLPYLERLDLALPGPFSPFPAISVRF PWYYEEYKKKLAAGGKPGELDKAFRSRDGRTAFEVLLMIARRLSQLGVKPRDGTEWSQNMPVGMI **TEEGIFPIPNLERFINAQLRRVRIVDEAGNVRAPTVEDIYKMGGYMVLVPTGRVEAVKDELWSKALG RDVVVRVHVYKPVKYSVDLEEWLWRTIHYNSPMAR**GEVPLPTPSGRVEIYSINLAYDVRRVFGKPA TSIDPSDLEGKKSGVDPLFSPVPLYAGMARPDYMWATGPATEDVEINGLVPPEPPKRLLLVYRHGPY THTHSNTQNNLLLDTLTSSELLSAWIHPDTAAALGVKDGDWIEVRPAAPKVAKQLESVGVKEVPTA RFRVRVTPMVRRDVVAIYHYWLVPRGRLRVKAWKLADVRAGYSDDNYLGPMLAGKLGTPGAMG NTVVEVSRVGGL

Solute binding protein (Pisl_0913):

1345.7	1345.8	0.1	YNVPWLSLILK (584-594)
1581.7	1581.8	0.1	YASFQEYIELYR (302-313)
2024.9	2025.0	0.1	DIGDAFASALEQLGFTVDR (203-221)
2153.0	2153.1	0.1	KDIGDAFASALEQLGFTVDR (202-221)
2318.0	2318.2	0.2	YSHMIISQGPFYLYAIDTAK (645-664)
2423.0	2423.1	0.1	MMYDPAVWNHPFTGEPMPFR (403-422)
2617.5	2618.5	1.0	AINVPTVVVGQPATISVSLEVPSGAGR (696-722)
3013.7	3013.6	0.1	AYKFWVLTYAENVPIVSEATQVFIPK (768-793)
3631.8	3629.8	2.0	GDPSVVLYTAAAGFNDIILNPAPSNPPCANPFSSR (70-104)
4383.5	4380.2	3.3	AAQAAPLKGDPSVVLYTAAAGFNDIILNPAPSNPPCANPFSSR (62-104)

Predicted amino acid sequence (matches are in red):

MRLRTVLLAMLLGIVLFAQYTPPHTNPGPATDRIVGKSVPIAQAGAAAKAGDIDVYIFGMRAAQAA PLKGDPSVVLYTAAAGFNDIILNPAPSNPPCANPFSSRAIRYAMQFVIDRDYVANEIFKGFAVPMYIW LSQYDPTYSIVADIISQLGIRYDLDYAKTIVESEMPKLGATKGPDGKWYCKGKLVTVIGLIRVEDERK DIGDAFASALEQLGFTVDRKYVTFDVAIQTVYGTDPAQFQWHFYTEGWGKSGIDRWDTSSIAQYCA SWFGYMPGWGTTGWYNFANATIDEITDKLYKGKYASFQEYIELYRKATLMCIQESVRVFVNTNLN AFVASPQLKGVTVDLGAGLRASVYNARNWYVPGKDVNVGHLWVWTASSAWNPVPQGGFTDVY SVDWFRMMYDPAVWNHPFTGEPMPFRATYVVETKGPAGYFDVPADAYRWDAKQKAWVSAGGA KAKSKVVFNYAKYIGAKWHHGQPIKLADVLFIYAFLWDIANDPQKVARESGVASYVNSTMNLIKGI RIINDTSIEVYIDYWHFDPNYIAAMAVTTPDMPWEVYYAVDQLVYVKQTYAASRASATKYNVPWL SLILKDHAKAAADVLQDALNKGIYPESWFKIGDKTLLTKDEALARYRAAVDWFNKYSHMIISQGPF YLYAIDTAKQYIELRAYRDPSYPYKPGAFYFGVATPVSVKAINVPTVVVGQPATISVSLEVPSGAGRI YYKWGIVDPTTGRFLYMSEEGTTAAAPININVPADVSSKLTANRAYKFWVLTYAENVPIVSEATQVF IPKAAAPATTPTPTPTPPPATTPTPPQPTVVTTTAPATGTTEALAAAIVGILAVLVALAFALRKKSGE TKQETKVYR