

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG031317\
 Data File : BG026223.D
 Acq On : 13 Mar 2017 22:25
 Operator : SJ/MA
 Sample : I2217-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Mar 15 17:17:17 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG030317-PAH.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 15 17:14:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	87190	20.00	ng/ul	0.00
2) Naphthalene-d8	10.86	136	429918	20.00	ng/ul	0.00
6) Acenaphthene-d10	14.66	164	261891	20.00	ng/ul	0.00
12) Phenanthrene-d10	17.39	188	600256	20.00	ng/ul	0.00
18) Chrysene-d12	21.64	240	629570	20.00	ng/ul	0.00
23) Perylene-d12	24.82	264	599259	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 2,4-Dichlorophenol-d3	10.48	165	8546	1.34	ng/uL	0.00
7) Acenaphthylene-d8	14.36	160	272177	12.19	ng/ul	0.00
10) Fluorene-d10	15.64	176	198910	12.21	ng/ul	0.00
15) Anthracene-d10	17.49	188	270863	10.15	ng/ul	0.00
19) Pyrene-d10	19.77	212	312404	9.72	ng/ul	0.00
26) Benzo(a)pyrene-d12	24.60	264	261022	10.10	ng/ul	0.00
Target Compounds				Qvalue		
14) Phenanthrene	17.44	178	236502	7.43	ng/ul	99
16) Anthracene	17.52	178	46608	1.45	ng/ul	97
17) Fluoranthene	19.43	202	389365	11.79	ng/ul	99
20) Pyrene	19.79	202	346585	8.65	ng/ul	99
21) Benzo(a)anthracene	21.62	228	188612	5.38	ng/ul	96
22) Chrysene	21.68	228	186677	5.71	ng/ul	98
24) Benzo(b)fluoranthene	23.82	252	233645	6.87	ng/ul	97
25) Benzo(k)fluoranthene	23.82	252	233645	7.19	ng/ul	96
27) Benzo(a)pyrene	24.67	252	152622	4.83	ng/ul	99
28) Indeno(1,2,3-cd)pyrene	28.45	276	99112	2.84	ng/ul#	91
30) Benzo(g,h,i)perylene	29.56	276	43607	1.46	ng/ul	93

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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