

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022117\
 Data File : BG025920.D
 Acq On : 22 Feb 2017 1:05
 Operator : SJ/MA
 Sample : I1901-08MSD
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DABF2MSD

Quant Time: Feb 22 04:49:45 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.3-EPA-BG021617MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Feb 21 23:44:02 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.09	152	152670	20.00	ng/ul	0.00
2) Naphthalene-d8	10.89	136	689432	20.00	ng/ul	0.00
6) Acenaphthene-d10	14.69	164	383260	20.00	ng/ul	0.00
12) Phenanthrene-d10	17.42	188	769061	20.00	ng/ul	0.00
17) Chrysene-d12	21.68	240	750788	20.00	ng/ul	0.00
22) Perylene-d12	24.93	264	683000	20.00	ng/ul	0.05
System Monitoring Compounds						
7) Acenaphthylene-d8	14.38	160	1219575	38.43	ng/ul	0.00
10) Fluorene-d10	15.67	176	869916	35.69	ng/ul	0.00
14) Anthracene-d10	17.52	188	1200714	34.53	ng/ul	0.00
18) Pyrene-d10	19.79	212	1251239	34.58	ng/ul	0.00
25) Benzo(a)pyrene-d12	24.72	264	1087416	35.98	ng/ul	0.06
Target Compounds						
9) Acenaphthene	14.75	153	880323	37.89	ng/ul	Ovalue 95
16) Fluoranthene	19.46	202	72617	1.61	ng/ul	98
19) Pyrene	19.83	202	1639669	36.09	ng/ul#	96
21) Chrysene	21.72	228	42623	1.06	ng/ul#	68

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

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022117\
Data File : BG025920.D
Acq On : 22 Feb 2017 1:05
Operator : SJ/MA
Sample : I1901-08MSD
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DABF2MSD

Quant Time: Feb 22 04:49:45 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.3-EPA-BG021617MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 21 23:44:02 2017
Response via : Initial Calibration

