

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072115\
 Data File : BG018048.D
 Acq On : 22 Jul 2015 7:55
 Operator : TP/UM
 Sample : G3027-06
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-05-3.5-4.0

Quant Time: Jul 22 08:28:28 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 22 04:14:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	59251	20.00	ng	0.00
21) Naphthalene-d8	10.33	136	283503	20.00	ng	0.00
38) Acenaphthene-d10	14.21	164	176200	20.00	ng	0.00
63) Phenanthrene-d10	16.97	188	349092	20.00	ng	0.00
75) Chrysene-d12	21.17	240	376003	20.00	ng	0.00
86) Perylene-d12	23.36	264	375657	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.16	112	509457	149.99	ng	0.00
7) Phenol-d6	6.74	99	725473	163.03	ng	0.00
23) Nitrobenzene-d5	8.70	82	435169	99.42	ng	0.00
41) 2,4,6-Tribromophenol	15.71	330	310474	173.92	ng	0.00
44) 2-Fluorobiphenyl	12.82	172	926195	91.09	ng	0.00
78) Terphenyl-d14	19.62	244	1213780	81.32	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	13.68	163	149105	12.18	ng	99
70) Phenanthrene	17.01	178	62399	3.47	ng	94
74) Fluoranthene	19.04	202	113644	5.76	ng	96
77) Pyrene	19.40	202	105161	5.01	ng	95
80) Benzo(a)anthracene	21.15	228	56833	2.86	ng	97
82) Chrysene	21.20	228	52723	2.79	ng	98
87) Benzo(b)fluoranthene	22.71	252	72543m	3.53	ng	
89) Benzo(a)pyrene	23.26	252	49067	2.56	ng	98

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

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