

Data Path : S:\HPCHEM1\BNA_G\DATA\BG072015\
 Data File : BG017987.D
 Acq On : 20 Jul 2015 14:05
 Operator : TP/UM
 Sample : G2976-09 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BRIDGE14-9(0-2)

Quant Time: Jul 20 18:18:23 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG071715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 18 01:50:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	151392	20.00	ng	0.00
21) Naphthalene-d8	10.37	136	651817	20.00	ng	0.00
38) Acenaphthene-d10	14.24	164	385511	20.00	ng	0.00
63) Phenanthrene-d10	17.00	188	733813	20.00	ng	0.00
75) Chrysene-d12	21.20	240	732166	20.00	ng	0.00
86) Perylene-d12	23.42	264	760495	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	145580	16.77	ng	0.00
7) Phenol-d6	6.77	99	194582	17.11	ng	0.00
23) Nitrobenzene-d5	8.74	82	111355	11.06	ng	0.00
41) 2,4,6-Tribromophenol	15.74	330	61484	15.74	ng	0.00
44) 2-Fluorobiphenyl	12.86	172	260592	11.71	ng	0.00
78) Terphenyl-d14	19.64	244	299959	10.32	ng	0.00
Target Compounds						
49) Dimethylphthalate	13.70	163	87752	3.28	ng	Qvalue 96
70) Phenanthrene	17.04	178	55823	1.47	ng	95
74) Fluoranthene	19.06	202	103360	2.49	ng	95
77) Pyrene	19.43	202	89305	2.19	ng	# 92
79) Butylbenzylphthalate	20.35	149	77776	4.25	ng	88
80) Benzo(a)anthracene	21.18	228	44027	1.14	ng	99
82) Chrysene	21.18	228	44027	1.20	ng	95
87) Benzo(b)fluoranthene	22.75	252	69882	1.68	ng	95
88) Benzo(k)fluoranthene	22.75	252	69882	1.68	ng	94
89) Benzo(a)pyrene	23.32	252	47344	1.22	ng	# 93

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

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