

Data Path : S:\HPCHEM1\BNA_G\DATA\BG072015\
 Data File : BG017983.D
 Acq On : 20 Jul 2015 11:43
 Operator : TP/UM
 Sample : G2976-05 2X
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BRIDGE14-7(0-2)

Quant Time: Jul 20 17:26:54 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	7.58	152	129848	20.00	ng	0.00
21) Naphthalene-d8	10.36	136	600573	20.00	ng	0.00
38) Acenaphthene-d10	14.24	164	370717	20.00	ng	0.00
63) Phenanthrene-d10	17.00	188	728424	20.00	ng	0.00
75) Chrysene-d12	21.20	240	731961	20.00	ng	0.00
86) Perylene-d12	23.42	264	735567	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	362148	48.65	ng	0.00
7) Phenol-d6	6.77	99	499613	51.23	ng	0.00
23) Nitrobenzene-d5	8.74	82	283540	30.58	ng	0.00
41) 2,4,6-Tribromophenol	15.74	330	164913	43.91	ng	0.00
44) 2-Fluorobiphenyl	12.86	172	629085	29.41	ng	0.00
78) Terphenyl-d14	19.64	244	771775	26.56	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	13.71	163	208119	8.08	ng	99
70) Phenanthrene	17.04	178	64170	1.71	ng	# 95
74) Fluoranthene	19.07	202	105922	2.58	ng	98
77) Pyrene	19.43	202	86448	2.12	ng	97
79) Butylbenzylphthalate	20.35	149	117528	6.43	ng	93
80) Benzo(a)anthracene	21.18	228	52235	1.35	ng	98
82) Chrysene	21.23	228	42270	1.15	ng	97
87) Benzo(b)fluoranthene	22.76	252	62872	1.56	ng	# 95
88) Benzo(k)fluoranthene	22.76	252	62872	1.56	ng	# 96
89) Benzo(a)pyrene	23.32	252	44458	1.19	ng	95

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

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