

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072115\
 Data File : BG018024.D
 Acq On : 21 Jul 2015 15:57
 Operator : TP/UM
 Sample : G2957-01RE
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SP-1RE

Quant Time: Jul 22 11:27:52 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 01:57:03 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	115844	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	519710	20.00	ng	0.00
38) Acenaphthene-d10	14.23	164	330034	20.00	ng	0.00
63) Phenanthrene-d10	16.98	188	643232	20.00	ng	0.00
75) Chrysene-d12	21.19	240	665085	20.00	ng	0.00
86) Perylene-d12	23.41	264	694922	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.17	112	716745	107.93	ng	0.00
7) Phenol-d6	6.76	99	940427	108.10	ng	0.00
23) Nitrobenzene-d5	8.73	82	527883	65.79	ng	0.00
41) 2,4,6-Tribromophenol	15.73	330	390549	116.80	ng	0.00
44) 2-Fluorobiphenyl	12.85	172	877497	46.08	ng	0.00
78) Terphenyl-d14	19.63	244	1125932	42.64	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.78	94	33265	3.58	ng	97
31) Naphthalene	10.40	128	57469	2.26	ng	98
49) Dimethylphthalate	13.69	163	282563	12.33	ng	99
57) Fluorene	15.29	166	60304	2.57	ng	98
70) Phenanthrene	17.03	178	489945	14.77	ng	99
71) Anthracene	17.12	178	133669	4.14	ng	99
74) Fluoranthene	19.06	202	815965	22.46	ng	97
77) Pyrene	19.42	202	705733	19.03	ng	98
80) Benzo(a)anthracene	21.17	228	460883	13.12	ng	99
82) Chrysene	21.23	228	400046	11.96	ng	96
85) Indeno(1,2,3-cd)pyrene	25.64	276	255882	6.32	ng	# 86
87) Benzo(b)fluoranthene	22.74	252	579385m	15.24	ng	
88) Benzo(k)fluoranthene	22.78	252	157366m	4.14	ng	
89) Benzo(a)pyrene	23.31	252	381697	10.78	ng	98
91) Benzo(g,h,i)perylene	26.34	276	247180	6.98	ng	97

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

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