

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG072115\
 Data File : BG018023.D
 Acq On : 21 Jul 2015 15:03
 Operator : TP/UM
 Sample : G2954-03 2X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TEC-E3

Quant Time: Jul 22 02:25:27 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	7.57	152	124848	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	520988	20.00	ng	0.00
38) Acenaphthene-d10	14.23	164	310682	20.00	ng	0.00
63) Phenanthrene-d10	16.99	188	601748	20.00	ng	0.00
75) Chrysene-d12	21.19	240	642692	20.00	ng	0.00
86) Perylene-d12	23.41	264	673521	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.18	112	331138	46.27	ng	0.00
7) Phenol-d6	6.76	99	438775	46.80	ng	0.00
23) Nitrobenzene-d5	8.72	82	255717	31.79	ng	0.00
41) 2,4,6-Tribromophenol	15.73	330	167557	53.23	ng	0.00
44) 2-Fluorobiphenyl	12.84	172	615934	34.36	ng	0.00
78) Terphenyl-d14	19.63	244	775079	30.38	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	13.70	163	141105	6.54	ng	99
70) Phenanthrene	17.03	178	450821	14.52	ng	99
71) Anthracene	17.11	178	114399	3.79	ng	99
74) Fluoranthene	19.05	202	879363	25.88	ng	98
77) Pyrene	19.42	202	866265	24.17	ng	98
80) Benzo(a)anthracene	21.17	228	485638	14.31	ng	98
82) Chrysene	21.23	228	456279	14.11	ng	96
85) Indeno(1,2,3-cd)pyrene	25.65	276	304905	7.80	ng	# 91
87) Benzo(b)fluoranthene	22.74	252	643175m	17.45	ng	
88) Benzo(k)fluoranthene	22.78	252	212910m	5.79	ng	
89) Benzo(a)pyrene	23.31	252	471853	13.74	ng	97
90) Dibenzo(a,h)anthracene	25.65	278	79899	2.24	ng	# 71
91) Benzo(g,h,i)perylene	26.33	276	316753	9.23	ng	97

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

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