

Data Path : Z:\HPCHEM1\BNA F\DATA\BF022718\
 Data File : BF103286.D
 Acq On : 28 Feb 2018 2:28
 Operator : SJ/JU
 Sample : J1683-02
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RFK-RMAB-MW08-GW

Quant Time: Feb 28 04:20:59 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 17:05:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	165068	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	624639	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	246509	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	362900	20.00	ng	0.00
75) Chrysene-d12	14.05	240	271334	20.00	ng	0.00
86) Perylene-d12	15.55	264	210135	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	287438	26.34	ng	0.00
7) Phenol-d6	6.52	99	483045	36.17	ng	0.00
23) Nitrobenzene-d5	7.45	82	863129	78.91	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	216540	103.78	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1159256	79.84	ng	0.00
78) Terphenyl-d14	12.99	244	915284	81.09	ng	0.00
Target Compounds						
10) Phenol	6.53	94	32046	2.067	ng	# 69
20) 3+4-Methylphenols	7.29	107	56150	4.471	ng	# 53
22) Acetophenone	7.26	105	481310	33.011	ng	# 58
27) 2,4-Dimethylphenol	7.83	122	271651	31.349	ng	# 91
31) Naphthalene	8.19	128	2683159	87.739	ng	# 98
37) 2-Methylnaphthalene	8.87	142	138011	6.983	ng	# 96

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

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