

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101717\
 Data File : BG029335.D
 Acq On : 18 Oct 2017 15:47
 Operator : SJ/JU
 Sample : I5906-16DL 5X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WH-TWS-06DL

Quant Time: Oct 18 16:54:43 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 17:32:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	65802	20.00	ng	0.00
21) Naphthalene-d8	10.98	136	323033	20.00	ng	0.00
38) Acenaphthene-d10	14.79	164	211100	20.00	ng	0.00
63) Phenanthrene-d10	17.52	188	467778	20.00	ng	0.00
75) Chrysene-d12	21.79	240	504146	20.00	ng	0.00
86) Perylene-d12	25.10	264	505984	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.72	112	102615	26.05	ng	0.00
7) Phenol-d6	7.32	99	136413	24.67	ng	0.00
23) Nitrobenzene-d5	9.33	82	90675	17.84	ng	0.00
41) 2,4,6-Tribromophenol	16.27	330	66919	24.55	ng	0.00
44) 2-Fluorobiphenyl	13.41	172	262348	18.23	ng	0.00
78) Terphenyl-d14	20.11	244	388101	17.51	ng	-0.01
Target Compounds						
10) Phenol	7.35	94	217520	36.492	ng	90
17) 2-Methylphenol	8.60	107	61113	15.434	ng	98
20) 3+4-Methylphenols	8.93	107	247771	44.409	ng	92
27) 2,4-Dimethylphenol	10.14	122	87810	17.767	ng	98
31) Naphthalene	11.04	128	793474	49.286	ng	98
37) 2-Methylnaphthalene	12.63	142	62691	5.343	ng	94
51) Acenaphthene	14.85	154	31111	2.308	ng	97
54) Dibenzofuran	15.18	168	65273	3.393	ng	97
70) Phenanthrene	17.56	178	114188	4.725	ng	99
72) Carbazole	17.92	167	182363	7.823	ng	99

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

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