

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092017\
 Data File : BG028849.D
 Acq On : 21 Sep 2017 9:07
 Operator : SJ/JU
 Sample : I5310-03
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NB-01-091917-A

Quant Time: Sep 21 10:33:41 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG091217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 12 16:57:11 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.29	152	33303	20.00	ng	-0.04
21) Naphthalene-d8	11.11	136	135976	20.00	ng	-0.04
38) Acenaphthene-d10	14.91	164	98663	20.00	ng	-0.04
63) Phenanthrene-d10	17.66	188	277096	20.00	ng	-0.04
75) Chrysene-d12	21.98	240	311462	20.00	ng	-0.05
86) Perylene-d12	25.45	264	303719	20.00	ng	-0.07
System Monitoring Compounds						
5) 2-Fluorophenol	5.86	112	283391	140.41	ng	-0.04
7) Phenol-d6	7.45	99	344797	113.46	ng	-0.04
23) Nitrobenzene-d5	9.46	82	279520	98.34	ng	-0.04
41) 2,4,6-Tribromophenol	16.40	330	227579	139.46	ng	-0.04
44) 2-Fluorobiphenyl	13.53	172	703822	95.12	ng	-0.04
78) Terphenyl-d14	20.25	244	1405328	96.22	ng	-0.03
Target Compounds						
13) 1,4-Dichlorobenzene	8.33	146	10028	4.025	ng	96
14) 1,2-Dichlorobenzene	8.65	146	5563	2.260	ng	98
30) 1,2,4-Trichlorobenzene	10.97	180	9704	3.491	ng	98
31) Naphthalene	11.16	128	104155	14.666	ng	95
37) 2-Methylnaphthalene	12.75	142	17271	3.257	ng	97
70) Phenanthrene	17.70	178	31985	2.257	ng	95
72) Carbazole	18.07	167	38577	2.992	ng	92

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

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