

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG102317\
 Data File : BG029406.D
 Acq On : 23 Oct 2017 19:32
 Operator : SJ/JU
 Sample : I6063-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S-1-A

Manual Integrations
 APPROVED

Sohil
 10/24/2017 8:47:43 PM

Quant Time: Oct 24 06:46:30 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	8.17	152	74582	20.00	ng	0.00
21) Naphthalene-d8	10.98	136	325131	20.00	ng	0.00
38) Acenaphthene-d10	14.78	164	214513	20.00	ng	0.00
63) Phenanthrene-d10	17.52	188	560979	20.00	ng	0.00
75) Chrysene-d12	21.79	240	630011	20.00	ng	-0.02
86) Perylene-d12	25.09	264	629635	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.72	112	519377	116.33	ng	0.00
7) Phenol-d6	7.32	99	626704	100.01	ng	0.00
23) Nitrobenzene-d5	9.33	82	435385	85.10	ng	0.00
41) 2,4,6-Tribromophenol	16.26	330	373176	134.74	ng	0.00
44) 2-Fluorobiphenyl	13.41	172	1128827	77.18	ng	0.00
78) Terphenyl-d14	20.11	244	2069342	74.72	ng	0.00
Target Compounds						
31) Naphthalene	11.03	128	130089	8.028	ng	98
37) 2-Methylnaphthalene	12.63	142	43093	3.649	ng	92
51) Acenaphthene	14.84	154	42484	3.102	ng	99
54) Dibenzofuran	15.18	168	68163	3.487	ng	98
57) Fluorene	15.82	166	39347	2.436	ng	96
70) Phenanthrene	17.56	178	102732	3.545	ng	100

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

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