Quantitation Report (QT Reviewed)

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA_G\DATA\BG062417\
Data File : BG027653.D
Acq On : 24 Jun 2017 10:52
Operator : SJ/JU
Sample : I3836-01DL 50X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : KY032MW006-170619DL

Quant Time: Jun 29 14:41:42 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG062117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 21 18:52:15 2017
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.51	152	17534	20.00	ng		0.00
21) Naphthalene-d8	11.33	136	86078	20.00	ng		0.00
38) Acenaphthene-d10	15.09	164	56444	20.00	ng		0.00
63) Phenanthrene-d10	17.84	188	146715	20.00	ng		0.00
75) Chrysene-d12	22.20	240	160750	20.00	ng		-0.01
86) Perylene-d12	25.84	264	154220	20.00	ng		0.00
System Monitoring Compounds							
5) 2-Fluorophenol	6.09	112	983	0.80	ng		0.00
7) Phenol-d6	7.65	99	4588	2.52	ng		0.00
23) Nitrobenzene-d5	9.67	82	2455	1.35	ng		0.00
41) 2,4,6-Tribromophenol	16.57	330	1537	1.83	ng		0.00
44) 2-Fluorobiphenyl	13.71	172	5824	1.41	ng		-0.01
78) Terphenyl-d14	20.41	244	8270	1.21	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	4.10	88	27108	56.673	ng		87
10) Phenol	7.67	94	7253	3.904	ng		98
11) bis(2-Chloroethyl)ether	7.91	93	9048	6.222	ng		86
17) 2-Methylphenol	8.93	107	6193	4.737	ng		93
20) 3+4-Methylphenols	9.25	107	61433	34.319	ng		90
27) 2,4-Dimethylphenol	10.49	122	14822	10.641	ng	#	90

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

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