

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG062417\
 Data File : BG027652.D
 Acq On : 24 Jun 2017 10:07
 Operator : SJ/JU
 Sample : I3836-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 KY032MW006-170619

Quant Time: Jun 29 14:39:24 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	28112	20.00	ng	0.00
21) Naphthalene-d8	11.34	136	111312	20.00	ng	0.01
38) Acenaphthene-d10	15.10	164	67818	20.00	ng	0.00
63) Phenanthrene-d10	17.84	188	157545	20.00	ng	0.00
75) Chrysene-d12	22.21	240	176413	20.00	ng	0.00
86) Perylene-d12	25.85	264	164189	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	6.12	112	84115	42.58	ng	0.03
7) Phenol-d6	7.73	99	102704	35.12	ng	0.08
23) Nitrobenzene-d5	9.70	82	172826	73.47	ng	0.02
41) 2,4,6-Tribromophenol	16.58	330	93548	92.52	ng	0.00
44) 2-Fluorobiphenyl	13.73	172	370383	74.60	ng	0.00
78) Terphenyl-d14	20.42	244	451260	60.27	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	4.10	88	1659927	2164.514	ng	# 77
10) Phenol	7.76	94	451953	151.721	ng	92
11) bis(2-Chloroethyl)ether	7.93	93	623397	267.375	ng	85
14) 1,2-Dichlorobenzene	8.87	146	51528	23.542	ng	94
17) 2-Methylphenol	8.98	107	395420	188.643	ng	96
20) 3+4-Methylphenols	9.34	107	3364619	1172.365	ng	96
27) 2,4-Dimethylphenol	10.55	122	929311	515.946	ng	93
31) Naphthalene	11.39	128	24066	4.000	ng	95
49) Dimethylphthalate	14.54	163	52757	8.885	ng	98
59) Diethylphthalate	15.88	149	31826	4.912	ng	98

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

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