

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122920\
 Data File : BG047839.D
 Acq On : 29 Dec 2020 16:58
 Operator : CG/JU
 Sample : L5227-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OWL-C1-GW

Quant Time: Dec 29 17:37:20 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG121720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 17 15:55:16 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.200	152	38100	20.00	ng	0.00
21) Naphthalene-d8	11.027	136	133672	20.00	ng	# 0.00
39) Acenaphthene-d10	14.822	164	101473	20.00	ng	0.00
64) Phenanthrene-d10	17.560	188	229511	20.00	ng	# 0.00
76) Chrysene-d12	21.832	240	223124	20.00	ng	# 0.00
86) Perylene-d12	25.181	264	247124	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.727	112	143193	65.02	ng	0.00
7) Phenol-d6	7.337	99	127948	39.74	ng	0.00
23) Nitrobenzene-d5	9.370	82	302045	83.11	ng	0.00
42) 2,4,6-Tribromophenol	16.303	330	207664	119.88	ng	0.00
45) 2-Fluorobiphenyl	13.447	172	611866	78.03	ng	0.00
79) Terphenyl-d14	20.145	244	933598	68.95	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	7.366	94	52195	17.88	ng	# 59
50) Dimethylphthalate	14.258	163	59477	6.80	ng	# 96
52) Acenaphthene	14.887	154	27954	4.24	ng	# 94
55) Dibenzofuran	15.216	168	28247	2.96	ng	# 70
58) Fluorene	15.862	166	52244	6.44	ng	95
71) Phenanthrene	17.601	178	78817	6.37	ng	97

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

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