

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041221\
 Data File : BG048669.D
 Acq On : 12 Apr 2021 20:35
 Operator : CG/JU
 Sample : M1966-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-115RR

Quant Time: Apr 13 00:20:00 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG033021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 07 10:06:00 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.088	152	23716	20.00	ng	#-0.01
21) Naphthalene-d8	10.919	136	94662	20.00	ng	# 0.00
39) Acenaphthene-d10	14.744	164	65708	20.00	ng	0.00
64) Phenanthrene-d10	17.500	188	147058	20.00	ng	0.00
76) Chrysene-d12	21.795	240	143586	20.00	ng	# 0.00
86) Perylene-d12	25.150	264	143997	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.643	112	109991	81.88	ng	0.00
7) Phenol-d6	7.247	99	92119	47.28	ng	0.00
23) Nitrobenzene-d5	9.263	82	173150	88.56	ng	0.00
42) 2,4,6-Tribromophenol	16.237	330	119119	137.91	ng	0.00
45) 2-Fluorobiphenyl	13.364	172	354028	80.08	ng	0.00
79) Terphenyl-d14	20.109	244	597575	76.11	ng	0.00
Target Compounds						
10) Phenol	7.277	94	7916	4.06	ng	Qvalue 95
27) 2,4-Dimethylphenol	10.079	122	22602	16.29	ng	96
50) Dimethylphthalate	14.186	163	45786	9.39	ng	# 97

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

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