

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG123120\
 Data File : BG047854.D
 Acq On : 31 Dec 2020 18:54
 Operator : CG/JU
 Sample : L5242-17
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SAMPLE-3(474+26)

Quant Time: Dec 31 23:49:25 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	37912	20.00	ng	# 0.00
21) Naphthalene-d8	11.020	136	142513	20.00	ng	# 0.00
39) Acenaphthene-d10	14.816	164	111798	20.00	ng	0.00
64) Phenanthrene-d10	17.554	188	271277	20.00	ng	# 0.00
76) Chrysene-d12	21.825	240	259740	20.00	ng	# 0.00
86) Perylene-d12	25.168	264	269510	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.732	112	244144	111.40	ng	0.00
7) Phenol-d6	7.342	99	306480	95.67	ng	0.00
23) Nitrobenzene-d5	9.369	82	307019	79.24	ng	0.00
42) 2,4,6-Tribromophenol	16.302	330	245936	128.86	ng	0.00
45) 2-Fluorobiphenyl	13.447	172	678833	78.57	ng	0.00
79) Terphenyl-d14	20.139	244	1234131	78.29	ng	0.00
Target Compounds						
10) Phenol	7.372	94	14717	5.07	ng	# 67
50) Dimethylphthalate	14.258	163	25752	2.67	ng	# 97
71) Phenanthrene	17.601	178	33160	2.27	ng	95

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

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