

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111120\
 Data File : BG047281.D
 Acq On : 10 Nov 2020 23:06
 Operator : CG/JU
 Sample : L4671-07
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 2A

Quant Time: Nov 11 02:01:40 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG110420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 04 16:19:40 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.011	152	67802	20.00	ng	0.00
21) Naphthalene-d8	10.826	136	266399	20.00	ng	0.01
39) Acenaphthene-d10	14.656	164	176982	20.00	ng	0.01
64) Phenanthrene-d10	17.424	188	398306	20.00	ng	# 0.04
76) Chrysene-d12	21.719	240	382108	20.00	ng	# 0.07
86) Perylene-d12	24.944	264	343855	20.00	ng	# 0.11
System Monitoring Compounds						
5) 2-Fluorophenol	5.579	112	276804	66.18	ng	0.02
7) Phenol-d6	7.189	99	382262	63.13	ng	0.03
23) Nitrobenzene-d5	9.181	82	271870	43.80	ng	0.01
42) 2,4,6-Tribromophenol	16.161	330	187322	63.21	ng	0.03
45) 2-Fluorobiphenyl	13.276	172	575482	41.70	ng	0.01
79) Terphenyl-d14	20.050	244	1055092	48.56	ng	0.05
Target Compounds						Qvalue
50) Dimethylphthalate	14.104	163	105628	6.76	ng	# 88
71) Phenanthrene	17.471	178	265182	11.47	ng	# 55
75) Fluoranthene	19.498	202	86719	2.95	ng	# 1

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

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