

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040821\
 Data File : BG048637.D
 Acq On : 8 Apr 2021 13:35
 Operator : CG/JU
 Sample : M1961-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TR1-A-D-COMP

Quant Time: Apr 08 14:09:09 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	8.094	152	22841	20.00	ng	0.00
21) Naphthalene-d8	10.914	136	81525	20.00	ng	-0.01
39) Acenaphthene-d10	14.745	164	53432	20.00	ng	0.00
64) Phenanthrene-d10	17.495	188	145436	20.00	ng	# 0.00
76) Chrysene-d12	21.790	240	159028	20.00	ng	#-0.01
86) Perylene-d12	25.133	264	150713	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.638	112	181068	139.96	ng	0.00
7) Phenol-d6	7.248	99	201883	107.60	ng	0.00
23) Nitrobenzene-d5	9.263	82	176213	104.65	ng	0.00
42) 2,4,6-Tribromophenol	16.237	330	142545	202.95	ng	0.00
45) 2-Fluorobiphenyl	13.364	172	384098	106.85	ng	0.00
79) Terphenyl-d14	20.109	244	814740	93.69	ng	0.00
Target Compounds						
10) Phenol	7.277	94	5210	2.77	ng	97
50) Dimethylphthalate	14.187	163	17446	4.40	ng	99

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

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