

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032221\
 Data File : BG048458.D
 Acq On : 23 Mar 2021 1:18
 Operator : CG/JU
 Sample : M1712-20
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RED-BRICK-COMP

Quant Time: Mar 23 02:08:26 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG031821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 22 11:22:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.103	152	56974	20.00	ng	# 0.00
21) Naphthalene-d8	10.923	136	230714	20.00	ng	# 0.00
39) Acenaphthene-d10	14.748	164	170545	20.00	ng	0.00
64) Phenanthrene-d10	17.498	188	370944	20.00	ng	# 0.00
76) Chrysene-d12	21.787	240	363809	20.00	ng	# 0.00
86) Perylene-d12	25.118	264	379029	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.659	112	435729	126.47	ng	0.01
7) Phenol-d6	7.263	99	541729	108.06	ng	0.01
23) Nitrobenzene-d5	9.272	82	487955	84.56	ng	0.00
42) 2,4,6-Tribromophenol	16.234	330	367851	132.63	ng	0.00
45) 2-Fluorobiphenyl	13.367	172	1049325	88.51	ng	0.00
79) Terphenyl-d14	20.106	244	1716946	85.11	ng	0.00
Target Compounds						
10) Phenol	7.286	94	10625	2.07	ng	# 56
50) Dimethylphthalate	14.184	163	34730	2.50	ng	99

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

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