

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072122\
 Data File : BG054349.D
 Acq On : 21 Jul 2022 17:46
 Operator : CG/JU
 Sample : N3767-03RE 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TANK-NRE

Quant Time: Jul 22 02:37:23 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG071522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 15 15:54:11 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.999	152	23206	20.000	ng	# 0.00
21) Naphthalene-d8	10.808	136	85543	20.000	ng	# 0.00
39) Acenaphthene-d10	14.633	164	68756	20.000	ng	0.00
64) Phenanthrene-d10	17.377	188	163071	20.000	ng	# 0.00
76) Chrysene-d12	21.648	240	183356	20.000	ng	# 0.00
86) Perylene-d12	24.856	264	193493	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.579	112	2680	2.411	ng	0.01
7) Phenol-d6	7.183	99	3239	2.127	ng	0.01
23) Nitrobenzene-d5	9.163	82	2391	1.522	ng	0.00
42) 2,4,6-Tribromophenol	16.119	330	3399	2.354	ng	0.00
45) 2-Fluorobiphenyl	13.252	172	8207	1.684	ng	0.00
79) Terphenyl-d14	19.985	244	17961	1.943	ng	0.00
Target Compounds						
32) Benzoic acid	10.162	122	32357	71.746	ng	# 84

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

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