

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090222\
 Data File : BG054841.D
 Acq On : 2 Sep 2022 15:30
 Operator : CG/JU
 Sample : N4453-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SASP2-T-5-(5-10)

Quant Time: Sep 03 00:27:29 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 25 17:50:55 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.147	152	32743	20.000	ng	0.00
21) Naphthalene-d8	10.973	136	135227	20.000	ng	# 0.00
39) Acenaphthene-d10	14.780	164	98538	20.000	ng	0.00
64) Phenanthrene-d10	17.530	188	239976	20.000	ng	0.00
76) Chrysene-d12	21.819	240	232710	20.000	ng	0.00
86) Perylene-d12	25.192	264	244562	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.685	112	198671	105.659	ng	0.00
7) Phenol-d6	7.295	99	275963	107.317	ng	0.00
23) Nitrobenzene-d5	9.322	82	166014	64.448	ng	0.00
42) 2,4,6-Tribromophenol	16.267	330	135658	110.175	ng	0.00
45) 2-Fluorobiphenyl	13.406	172	438943	66.953	ng	0.00
79) Terphenyl-d14	20.127	244	858071	73.060	ng	0.00

Target Compounds	Qvalue

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

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