

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090122\
 Data File : BG054820.D
 Acq On : 1 Sep 2022 13:23
 Operator : CG/JU
 Sample : N4431-01RE
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CRUSHED-MASONARY-PILE-4RE

Quant Time: Sep 01 14:34:54 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	41688	20.000	ng	0.00
21) Naphthalene-d8	10.973	136	168987	20.000	ng	0.00
39) Acenaphthene-d10	14.780	164	117316	20.000	ng	0.00
64) Phenanthrene-d10	17.524	188	270190	20.000	ng	0.00
76) Chrysene-d12	21.819	240	256900	20.000	ng	0.00
86) Perylene-d12	25.185	264	275367	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.691	112	23474	9.805	ng	0.00
7) Phenol-d6	7.295	99	121873	37.225	ng	0.00
23) Nitrobenzene-d5	9.322	82	200722	62.355	ng	0.00
42) 2,4,6-Tribromophenol	16.266	330	12125	8.271	ng	0.00
45) 2-Fluorobiphenyl	13.405	172	512803	65.699	ng	0.00
79) Terphenyl-d14	20.127	244	892050	68.801	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.571	178	54595	3.793	ng	96
75) Fluoranthene	19.574	202	75620	4.319	ng	96
78) Pyrene	19.933	202	74592	4.164	ng	98
81) Benzo(a)anthracene	21.795	228	44582	2.559	ng	97
83) Chrysene	21.866	228	38147	2.290	ng	95
88) Benzo(b)fluoranthene	24.104	252	36341	2.104	ng	99

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

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