

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041822\
 Data File : BG053190.D
 Acq On : 18 Apr 2022 21:30
 Operator : CG/JU
 Sample : N2444-08
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-7

Quant Time: Apr 19 04:21:46 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 08 15:31:24 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.120	152	29774	20.000	ng	0.00
21) Naphthalene-d8	10.928	136	116435	20.000	ng	0.00
39) Acenaphthene-d10	14.735	164	83367	20.000	ng	-0.01
64) Phenanthrene-d10	17.484	188	212275	20.000	ng	0.00
76) Chrysene-d12	21.760	240	245145	20.000	ng	-0.01
86) Perylene-d12	25.062	264	254617	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.682	112	241022	138.764	ng	0.00
7) Phenol-d6	7.274	99	343156	128.114	ng	0.00
23) Nitrobenzene-d5	9.277	82	243786	85.014	ng	0.00
42) 2,4,6-Tribromophenol	16.227	330	167839	126.543	ng	0.00
45) 2-Fluorobiphenyl	13.360	172	494411	81.776	ng	0.00
79) Terphenyl-d14	20.086	244	1025522	70.652	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	17.525	178	57045	4.920	ng	96
75) Fluoranthene	19.528	202	58505	3.936	ng	98
78) Pyrene	19.892	202	50219	2.876	ng	96

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

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