

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG052822\
 Data File : BG053759.D
 Acq On : 28 May 2022 20:31
 Operator : CG/JU
 Sample : N3056-02DL 20X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-12DL

Quant Time: May 29 02:20:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG050522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun May 29 02:17:06 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.949	152	53434	20.000	ng	0.00
21) Naphthalene-d8	10.757	136	209929	20.000	ng	# 0.00
39) Acenaphthene-d10	14.593	164	147843	20.000	ng	0.00
64) Phenanthrene-d10	17.337	188	315285	20.000	ng	# 0.00
76) Chrysene-d12	21.601	240	292685	20.000	ng	# 0.00
86) Perylene-d12	24.779	264	302177	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.535	112	7849	2.493	ng	0.00
7) Phenol-d6	7.162	99	6724	1.407	ng	0.00
23) Nitrobenzene-d5	9.118	82	18674	3.787	ng	0.00
42) 2,4,6-Tribromophenol	16.091	330	11366	5.213	ng	0.00
45) 2-Fluorobiphenyl	13.213	172	44389	4.387	ng	0.00
79) Terphenyl-d14	19.945	244	59151	3.764	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	10.810	128	443535	41.535	ng	100
37) 2-Methylnaphthalene	12.420	142	51150	6.418	ng	97
38) 1-Methylnaphthalene	12.637	142	148434	19.091	ng	98
46) 1,1'-Biphenyl	13.424	154	26164	2.596	ng	97
52) Acenaphthene	14.652	154	26184	3.487	ng	96
71) Phenanthrene	17.384	178	33821	2.110	ng	97

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

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