

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121823\
 Data File : BG060076.D
 Acq On : 18 Dec 2023 21:37
 Operator : MA/JU
 Sample : 05612-05DL 10X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-06SDL

Quant Time: Dec 19 04:42:18 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 19 04:37:23 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.953	152	64850	20.000	ng	0.00
21) Naphthalene-d8	10.762	136	246061	20.000	ng	# 0.00
39) Acenaphthene-d10	14.593	164	212813	20.000	ng	0.00
64) Phenanthrene-d10	17.337	188	651402	20.000	ng	#-0.01
76) Chrysene-d12	21.608	240	693769	20.000	ng	#-0.01
86) Perylene-d12	24.804	264	791229	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.521	112	18589	4.907	ng	0.00
7) Phenol-d6	7.107	99	17464	3.277	ng	0.00
23) Nitrobenzene-d5	9.105	82	40331	7.443	ng	-0.02
42) 2,4,6-Tribromophenol	16.079	330	64671	13.040	ng	0.00
45) 2-Fluorobiphenyl	13.212	172	154618	8.892	ng	0.00
79) Terphenyl-d14	19.957	244	377090	9.923	ng	-0.01
Target Compounds						
						Qvalue
31) Naphthalene	10.821	128	3253246	253.973	ng	# 96
37) 2-Methylnaphthalene	12.413	142	233300	23.222	ng	97
38) 1-Methylnaphthalene	12.636	142	399997	40.306	ng	96
46) 1,1'-Biphenyl	13.423	154	37416	2.371	ng	98
52) Acenaphthene	14.657	154	231102	17.957	ng	97
58) Fluorene	15.633	166	71389	4.042	ng	96
71) Phenanthrene	17.384	178	105150	3.333	ng	99

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

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