

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121823\
 Data File : BG060074.D
 Acq On : 18 Dec 2023 20:15
 Operator : MA/JU
 Sample : 05612-04DL 10X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-04IDL

Quant Time: Dec 19 04:41:30 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	7.958	152	64208	20.000	ng	0.00
21) Naphthalene-d8	10.767	136	254217	20.000	ng	# 0.00
39) Acenaphthene-d10	14.592	164	222299	20.000	ng	0.00
64) Phenanthrene-d10	17.342	188	661589	20.000	ng	# 0.00
76) Chrysene-d12	21.613	240	693555	20.000	ng	# 0.00
86) Perylene-d12	24.803	264	767305	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.526	112	24027	6.406	ng	0.00
7) Phenol-d6	7.106	99	21489	4.073	ng	0.00
23) Nitrobenzene-d5	9.116	82	48135	8.598	ng	0.00
42) 2,4,6-Tribromophenol	16.078	330	79438	15.334	ng	0.00
45) 2-Fluorobiphenyl	13.217	172	178263	9.815	ng	0.00
79) Terphenyl-d14	19.962	244	399086	10.506	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	7.136	94	20030	3.779	ng	77
27) 2,4-Dimethylphenol	9.927	122	13100	4.572	ng	# 84
31) Naphthalene	10.832	128	6389538	482.812	ng	# 85
37) 2-Methylnaphthalene	12.418	142	477845	46.037	ng	98
38) 1-Methylnaphthalene	12.635	142	390715	38.107	ng	97
46) 1,1'-Biphenyl	13.429	154	54970	3.334	ng	99
52) Acenaphthene	14.656	154	193081	14.362	ng	86
58) Fluorene	15.638	166	59914	3.247	ng	89
71) Phenanthrene	17.383	178	129367	4.037	ng	96

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

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