

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062323\
 Data File : BG057863.D
 Acq On : 24 Jun 2023 5:46
 Operator : CG/JU
 Sample : 03291-03DL 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MW-6ADL

Quant Time: Jun 25 01:43:27 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 20 03:40:38 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.924	152	50896	20.000	ng	0.00
21) Naphthalene-d8	10.727	136	215573	20.000	ng	0.00
39) Acenaphthene-d10	14.557	164	138629	20.000	ng	0.00
64) Phenanthrene-d10	17.301	188	300741	20.000	ng	0.00
76) Chrysene-d12	21.555	240	254062	20.000	ng	0.00
86) Perylene-d12	24.710	264	311592	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.503	112	112160	33.979	ng	0.00
7) Phenol-d6	7.084	99	94182	18.700	ng	0.00
23) Nitrobenzene-d5	9.082	82	199143	47.412	ng	-0.01
42) 2,4,6-Tribromophenol	16.044	330	142099	73.585	ng	0.00
45) 2-Fluorobiphenyl	13.183	172	441924	48.351	ng	0.00
79) Terphenyl-d14	19.916	244	636081	46.685	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	7.113	94	13283	2.679	ng	94
15) Benzyl Alcohol	8.153	79	9392	2.439	ng	96
17) 2-Methylphenol	8.365	107	151686	41.588	ng	95
20) 3+4-Methylphenols	8.688	107	33839	6.717	ng	96
31) Naphthalene	10.780	128	807771	70.404	ng	99
37) 2-Methylnaphthalene	12.384	142	70865	8.474	ng	99
38) 1-Methylnaphthalene	12.601	142	34495	4.370	ng	98

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

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