

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020924\
 Data File : BP019638.D
 Acq On : 09 Feb 2024 18:26
 Operator : MA/JU
 Sample : P1372-06DL 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CG-MW-17-ch5DL

Quant Time: Feb 10 00:51:40 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 01 00:51:48 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	83735	20.000	ng	-0.01
21) Naphthalene-d8	10.710	136	313065	20.000	ng	0.00
39) Acenaphthene-d10	14.539	164	201018	20.000	ng	-0.01
64) Phenanthrene-d10	17.298	188	432862	20.000	ng	0.00
76) Chrysene-d12	21.386	240	430274	20.000	ng	0.00
86) Perylene-d12	23.810	264	522931	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	73006	14.054	ng	0.00
7) Phenol-d6	7.075	99	57040	8.143	ng	-0.01
23) Nitrobenzene-d5	9.075	82	154404	21.371	ng	-0.01
42) 2,4,6-Tribromophenol	16.034	330	105248	35.860	ng	0.00
45) 2-Fluorobiphenyl	13.169	172	349845	21.515	ng	-0.01
79) Terphenyl-d14	19.863	244	542296	20.727	ng	0.00
Target Compounds						
						Qvalue
10) Phenol	7.105	94	15981	2.289	ng	97
20) 3+4-Methylphenols	8.675	107	322752	50.317	ng	91
31) Naphthalene	10.757	128	904851	52.699	ng	99
38) 1-Methylnaphthalene	12.587	142	136553	11.694	ng	97
49) Acenaphthylene	14.263	152	78530	4.250	ng	98
52) Acenaphthene	14.604	154	47108	4.108	ng	98

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

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