

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020924\
 Data File : BP019639.D
 Acq On : 09 Feb 2024 19:00
 Operator : MA/JU
 Sample : P1372-07DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CG-MW-17-ch6DL

Quant Time: Feb 10 00:52:10 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	7.904	152	77784	20.000	ng	-0.01
21) Naphthalene-d8	10.710	136	288851	20.000	ng	0.00
39) Acenaphthene-d10	14.539	164	183953	20.000	ng	-0.01
64) Phenanthrene-d10	17.298	188	433311	20.000	ng	0.00
76) Chrysene-d12	21.386	240	461314	20.000	ng	0.00
86) Perylene-d12	23.804	264	534316	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	61046	12.651	ng	0.00
7) Phenol-d6	7.075	99	45916	7.057	ng	-0.01
23) Nitrobenzene-d5	9.075	82	139968	20.997	ng	-0.01
42) 2,4,6-Tribromophenol	16.033	330	104227	38.806	ng	0.00
45) 2-Fluorobiphenyl	13.169	172	299639	20.137	ng	-0.01
79) Terphenyl-d14	19.863	244	585201	20.862	ng	0.00
Target Compounds						Qvalue
20) 3+4-Methylphenols	8.675	107	229145	38.457	ng	92
31) Naphthalene	10.757	128	957523	60.441	ng	99
38) 1-Methylnaphthalene	12.587	142	47021	4.364	ng	95

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

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