

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020924\
 Data File : BP019640.D
 Acq On : 09 Feb 2024 19:34
 Operator : MA/JU
 Sample : P1372-08DL 10X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CG-MW-17-ch7DL

Quant Time: Feb 10 00:52:38 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	81398	20.000	ng	-0.01
21) Naphthalene-d8	10.704	136	332444	20.000	ng	-0.01
39) Acenaphthene-d10	14.539	164	226192	20.000	ng	-0.01
64) Phenanthrene-d10	17.298	188	466366	20.000	ng	0.00
76) Chrysene-d12	21.386	240	377182	20.000	ng	0.00
86) Perylene-d12	23.810	264	449035	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	29372	5.817	ng	-0.01
7) Phenol-d6	7.075	99	25619	3.763	ng	-0.01
23) Nitrobenzene-d5	9.075	82	65853	8.583	ng	-0.01
42) 2,4,6-Tribromophenol	16.033	330	47411	14.356	ng	0.00
45) 2-Fluorobiphenyl	13.169	172	156392	8.547	ng	-0.01
79) Terphenyl-d14	19.863	244	195307	8.516	ng	0.00
Target Compounds						
20) 3+4-Methylphenols	8.675	107	451708	72.444	ng	91
31) Naphthalene	10.757	128	249270	13.671	ng	99

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

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