

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101220\
 Data File : BN012150.D
 Acq On : 12 Oct 2020 16:47
 Operator : CG/JU
 Sample : L3971-08DL2 100X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-ACIDS-WPDL2

Quant Time: Oct 12 17:23:51 2020
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN092320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 12 11:15:56 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.627	152	1557	0.40	ng	# 0.00
7) Naphthalene-d8	10.402	136	6201	0.40	ng	# 0.00
13) Acenaphthene-d10	14.268	164	3279	0.40	ng	0.00
19) Phenanthrene-d10	17.017	188	6521	0.40	ng	# 0.00
29) Chrysene-d12	21.227	240	6484	0.40	ng	# 0.00
36) Perylene-d12	23.426	264	7355	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.235	112	6878	1.26	ng	0.00
5) Phenol-d6	6.805	99	8180	1.19	ng	0.00
8) Nitrobenzene-d5	8.780	82	6204	0.97	ng	0.00
11) 2-Methylnaphthalene-d10	11.998	152	2	0.00	ng	0.00
14) 2,4,6-Tribromophenol	15.765	330	2424	1.52	ng	0.00
15) 2-Fluorobiphenyl	12.885	172	11399	0.92	ng	0.00
27) Fluoranthene-d10	19.072	212	7	0.00	ng	0.02
31) Terphenyl-d14	19.668	244	14521	0.98	ng	0.00
Target Compounds						
20) 4,6-Dinitro-2-methylph...	15.410	198	2027	1.63	ng	# 85
24) Pentachlorophenol	16.676	266	1114	0.57	ng	95

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101220\
Data File : BN012150.D
Acq On : 12 Oct 2020 16:47
Operator : CG/JU
Sample : L3971-08DL2 100X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
PT-ACIDS-WPDL2

Quant Time: Oct 12 17:23:51 2020
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN092320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 12 11:15:56 2020
Response via : Initial Calibration

