

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091820\
 Data File : BN011837.D
 Acq On : 18 Sep 2020 15:57
 Operator : CG/JU
 Sample : L3971-08DL 20X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Sep 18 16:26:36 2020
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN091720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 18 10:47:35 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.715	152	1959	0.40	ng	0.00
7) Naphthalene-d8	10.491	136	7507	0.40	ng	# 0.00
13) Acenaphthene-d10	14.357	164	4371	0.40	ng	0.00
19) Phenanthrene-d10	17.102	188	10201	0.40	ng	0.00
29) Chrysene-d12	21.312	240	9877	0.40	ng	# 0.00
36) Perylene-d12	23.566	264	9930	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.307	112	36537	6.31	ng	0.00
5) Phenol-d6	6.877	99	45333	6.39	ng	0.00
8) Nitrobenzene-d5	8.856	82	35580	4.70	ng	0.00
11) 2-Methylnaphthalene-d10	0.000	152	0d	0.00	ng	
14) 2,4,6-Tribromophenol	15.854	330	16345	7.35	ng	0.00
15) 2-Fluorobiphenyl	12.974	172	87078	4.61	ng	0.00
27) Fluoranthene-d10	0.000	212	0d	0.00	ng	
31) Terphenyl-d14	19.752	244	120880	5.36	ng	0.00
Target Compounds						
20) 4,6-Dinitro-2-methylph...	15.486	198	14581	8.35	ng	# 68
24) Pentachlorophenol	16.761	266	8267	2.73	ng	98

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

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