

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100821\
 Data File : BN016859.D
 Acq On : 09 Oct 2021 00:11
 Operator : CG/JU
 Sample : M4096-09
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MWHR2-6-100421

Quant Time: Oct 09 00:46:09 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 08 15:46:07 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.615	152	19176	0.400	ng	0.00
7) Naphthalene-d8	10.381	136	87814	0.400	ng	# 0.00
13) Acenaphthene-d10	14.243	164	47494	0.400	ng	0.00
19) Phenanthrene-d10	16.993	188	95414	0.400	ng	#-0.01
29) Chrysene-d12	21.206	240	98817	0.400	ng	# 0.00
35) Perylene-d12	23.447	264	105021	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.245	112	6246	0.130	ng	0.00
5) Phenol-d6	6.803	99	4102	0.075	ng	0.00
8) Nitrobenzene-d5	8.759	82	16498	0.268	ng	0.00
11) 2-Methylnaphthalene-d10	11.976	152	41369	0.316	ng	0.00
14) 2,4,6-Tribromophenol	15.741	330	9052	0.368	ng	-0.01
15) 2-Fluorobiphenyl	12.873	172	54422	0.284	ng	0.00
27) Fluoranthene-d10	19.043	212	111917	0.422	ng	0.00
31) Terphenyl-d14	19.654	244	92284	0.428	ng	0.00
Target Compounds						
24) Pentachlorophenol	16.653	266	120	0.107	ng	94
34) Bis(2-ethylhexyl)phtha...	21.142	149	21406	0.138	ng	99

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

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