

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101321\
 Data File : BN016917.D
 Acq On : 13 Oct 2021 03:13
 Operator : CG/JU
 Sample : M4169-05
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SW-RR-01-(2.0)

Quant Time: Oct 13 06:19:55 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN101321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 13 02:19:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.688	152	21336	0.400	ng	0.00
7) Naphthalene-d8	10.445	136	96381	0.400	ng	#-0.01
13) Acenaphthene-d10	14.294	164	51331	0.400	ng	0.00
19) Phenanthrene-d10	17.030	188	99394	0.400	ng	#-0.01
29) Chrysene-d12	21.227	240	99347	0.400	ng	# 0.00
35) Perylene-d12	23.467	264	103231	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.310	112	7104	0.147	ng	0.02
5) Phenol-d6	6.877	99	4464	0.083	ng	0.00
8) Nitrobenzene-d5	8.822	82	18010	0.294	ng	0.00
11) 2-Methylnaphthalene-d10	12.038	152	45231	0.335	ng	0.00
14) 2,4,6-Tribromophenol	15.791	330	8988	0.386	ng	0.00
15) 2-Fluorobiphenyl	12.924	172	60421	0.311	ng	0.00
27) Fluoranthene-d10	19.068	212	111916	0.448	ng	0.00
31) Terphenyl-d14	19.678	244	96853	0.444	ng	0.00
Target Compounds						
24) Pentachlorophenol	16.689	266	159	0.203	ng	# 86
34) Bis(2-ethylhexyl)phtha...	21.164	149	21044	0.302	ng	100

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

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