

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100121\
 Data File : BN016684.D
 Acq On : 02 Oct 2021 23:09
 Operator : CG/JU
 Sample : M3892-10
 Misc :
 ALS Vial : 52 Sample Multiplier: 2

Instrument :
 BNA_N
 ClientSampleId :
 S-824-N-SO-1-1.5-092221

Quant Time: Oct 04 03:35:57 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 01 13:35:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.642	152	62288	0.800	ng	0.00
7) Naphthalene-d8	10.407	136	288843	0.800	ng	# 0.00
13) Acenaphthene-d10	14.269	164	165212	0.800	ng	0.00
19) Phenanthrene-d10	17.018	188	323704	0.800	ng	#-0.01
29) Chrysene-d12	21.228	240	318574	0.800	ng	#-0.01
35) Perylene-d12	23.485	264	336649	0.800	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.273	112	22076	0.278	ng	0.02
5) Phenol-d6	6.831	99	22729	0.262	ng	0.00
8) Nitrobenzene-d5	8.785	82	25153	0.267	ng	0.00
11) 2-Methylnaphthalene-d10	12.003	152	60333	0.282	ng	0.00
14) 2,4,6-Tribromophenol	15.766	330	12053	0.320	ng	0.00
15) 2-Fluorobiphenyl	12.886	172	75383	0.225	ng	-0.01
27) Fluoranthene-d10	19.063	212	114452	0.266	ng	0.00
31) Terphenyl-d14	19.679	244	87397	0.252	ng	0.00
Target Compounds						
						Qvalue
32) Benzo(a)anthracene	21.217	228	19022	0.039	ng	98
33) Chrysene	21.270	228	18393	0.037	ng	98
34) Bis(2-ethylhexyl)phtha...	21.164	149	35425	0.182	ng	99
36) Indeno(1,2,3-cd)pyrene	25.764	276	18689	0.031	ng	# 84
37) Benzo(b)fluoranthene	22.807	252	20496	0.038	ng	# 70
38) Benzo(k)fluoranthene	22.851	252	18366	0.034	ng	# 35
40) Dibenzo(a,h)anthracene	25.788	278	14644	0.030	ng	# 70
41) Benzo(g,h,i)perylene	26.459	276	14224	0.027	ng	# 87

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

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