

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121821\
 Data File : BM033524.D
 Acq On : 18 Dec 2021 18:14
 Operator : CG/JU
 Sample : M5114-03
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 HS-OILY-WATER

Quant Time: Dec 20 07:08:28 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM121621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 16 17:10:15 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.931	152	29309	20.000	ng	-0.01
21) Naphthalene-d8	10.742	136	123361	20.000	ng	-0.01
39) Acenaphthene-d10	14.577	164	93168	20.000	ng	-0.01
64) Phenanthrene-d10	17.324	188	215184	20.000	ng	#-0.01
76) Chrysene-d12	21.500	240	211610	20.000	ng	#-0.01
86) Perylene-d12	23.912	264	190131	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.495	112	142744	84.039	ng	0.00
7) Phenol-d6	7.101	99	117868	51.709	ng	-0.01
23) Nitrobenzene-d5	9.101	82	318107	110.722	ng	-0.01
42) 2,4,6-Tribromophenol	16.071	330	187405	178.815	ng	0.00
45) 2-Fluorobiphenyl	13.201	172	671413	99.688	ng	-0.01
79) Terphenyl-d14	19.947	244	938004	69.785	ng	0.00

Target Compounds	Qvalue

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

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