

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102922\
 Data File : BM037267.D
 Acq On : 29 Oct 2022 12:53
 Operator : CG/JU
 Sample : N5266-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 20221100-204.1-SITE-WATER

Quant Time: Oct 31 02:05:18 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM102822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 31 01:34:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	500457	20.000	ng	0.00
21) Naphthalene-d8	10.675	136	1964054	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	1114217	20.000	ng	0.00
64) Phenanthrene-d10	17.245	188	2114079	20.000	ng	0.00
76) Chrysene-d12	21.415	240	1752172	20.000	ng	0.00
86) Perylene-d12	23.774	264	1855354	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.445	112	2292106	79.126	ng	0.00
7) Phenol-d6	7.045	99	2285852	58.141	ng	0.00
23) Nitrobenzene-d5	9.045	82	3641164	93.534	ng	0.00
42) 2,4,6-Tribromophenol	15.998	330	2095172	159.563	ng	0.00
45) 2-Fluorobiphenyl	13.133	172	8097469	96.562	ng	0.00
79) Terphenyl-d14	19.868	244	10318613	105.607	ng	0.00
Target Compounds						
10) Phenol	7.069	94	103435	2.348	ng	Qvalue 94

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

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