

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102922\
 Data File : BM037268.D
 Acq On : 29 Oct 2022 13:29
 Operator : CG/JU
 Sample : N5266-04
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 20221092-FIELD-BLANK-2

Quant Time: Oct 31 02:05:30 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	455593	20.000	ng	0.00
21) Naphthalene-d8	10.675	136	1761079	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	977854	20.000	ng	0.00
64) Phenanthrene-d10	17.245	188	1860114	20.000	ng	0.00
76) Chrysene-d12	21.421	240	1589755	20.000	ng	0.00
86) Perylene-d12	23.780	264	1687047	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	1793112	67.996	ng	0.00
7) Phenol-d6	7.046	99	1301189	36.355	ng	0.00
23) Nitrobenzene-d5	9.045	82	3377242	96.754	ng	0.00
42) 2,4,6-Tribromophenol	15.998	330	1977785	171.628	ng	0.00
45) 2-Fluorobiphenyl	13.133	172	7438219	101.069	ng	0.00
79) Terphenyl-d14	19.868	244	9766511	110.168	ng	0.00

Target Compounds	Qvalue

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

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