

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092322\
 Data File : BP011789.D
 Acq On : 23 Sep 2022 18:45
 Operator : CG/JU
 Sample : N4732-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 WC-1

Quant Time: Sep 24 01:37:22 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 22 02:45:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.916	152	520645	20.000	ng	0.00
21) Naphthalene-d8	10.728	136	2090094	20.000	ng	0.00
39) Acenaphthene-d10	14.557	164	1038050	20.000	ng	-0.01
64) Phenanthrene-d10	17.316	188	2061412	20.000	ng	0.00
76) Chrysene-d12	21.398	240	1504534	20.000	ng	0.00
86) Perylene-d12	23.845	264	224215	20.000	ng	# 0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	4923930	165.939	ng	0.00
7) Phenol-d6	7.081	99	5905232	134.593	ng	0.00
23) Nitrobenzene-d5	9.087	82	4202915	105.250	ng	0.00
42) 2,4,6-Tribromophenol	16.057	330	1340669	191.807	ng	0.00
45) 2-Fluorobiphenyl	13.187	172	7463544	113.701	ng	0.00
79) Terphenyl-d14	19.875	244	9047061	138.264	ng	0.00

Target Compounds	Qvalue

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

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