

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091922\
 Data File : BP011744.D
 Acq On : 19 Sep 2022 20:24
 Operator : CG/JU
 Sample : N4630-17
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 WC-17-4-4

Quant Time: Sep 20 04:49:01 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 20 04:42:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	504469	20.000	ng	0.00
21) Naphthalene-d8	10.746	136	2103676	20.000	ng	# 0.00
39) Acenaphthene-d10	14.575	164	1386615	20.000	ng	0.00
64) Phenanthrene-d10	17.328	188	3100505	20.000	ng	0.00
76) Chrysene-d12	21.404	240	3255931	20.000	ng	0.00
86) Perylene-d12	23.857	264	3378920	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	4402684	158.856	ng	0.00
7) Phenol-d6	7.099	99	5302211	143.519	ng	0.00
23) Nitrobenzene-d5	9.105	82	3465190	97.691	ng	0.00
42) 2,4,6-Tribromophenol	16.069	330	2959807	273.193	ng	0.00
45) 2-Fluorobiphenyl	13.204	172	8550545	95.854	ng	0.00
79) Terphenyl-d14	19.875	244	13401747	89.508	ng	0.00

Target Compounds	Qvalue

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

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