

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091922\
 Data File : BP011755.D
 Acq On : 20 Sep 2022 02:47
 Operator : CG/JU
 Sample : N4630-12
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 WC-17-4-3

Quant Time: Sep 20 04:55:15 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	7.934	152	483352	20.000	ng	0.00
21) Naphthalene-d8	10.746	136	2116382	20.000	ng	# 0.00
39) Acenaphthene-d10	14.575	164	453905	20.000	ng	# 0.00
64) Phenanthrene-d10	17.328	188	2363972	20.000	ng	# 0.00
76) Chrysene-d12	21.404	240	1203589	20.000	ng	# 0.00
86) Perylene-d12	23.851	264	16489	20.000	ng	# 0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	4603248	173.349	ng	0.00
7) Phenol-d6	7.099	99	3314334	93.631	ng	0.00
23) Nitrobenzene-d5	9.104	82	3842881	107.689	ng	0.00
42) 2,4,6-Tribromophenol	16.063	330	1968198	554.965	ng	0.00
45) 2-Fluorobiphenyl	13.198	172	8251892	282.593	ng	0.00
79) Terphenyl-d14	19.880	244	10732210	193.904	ng	0.00
Target Compounds						
32) Benzoic acid	9.963	122	29971	7.800	ng	Qvalue 96

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

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