

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091922\
 Data File : BP011746.D
 Acq On : 19 Sep 2022 21:34
 Operator : CG/JU
 Sample : N4697-04
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SB-23-COMP

Quant Time: Sep 20 04:49:57 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	385330	20.000	ng	0.00
21) Naphthalene-d8	10.746	136	1712214	20.000	ng	# 0.00
39) Acenaphthene-d10	14.575	164	1131335	20.000	ng	0.00
64) Phenanthrene-d10	17.328	188	2441631	20.000	ng	# 0.00
76) Chrysene-d12	21.404	240	2337389	20.000	ng	# 0.00
86) Perylene-d12	23.851	264	2303133	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	3862846	182.471	ng	0.00
7) Phenol-d6	7.099	99	5239738	185.680	ng	0.00
23) Nitrobenzene-d5	9.105	82	2844530	98.528	ng	0.00
42) 2,4,6-Tribromophenol	16.069	330	1762452	199.383	ng	0.00
45) 2-Fluorobiphenyl	13.204	172	6721322	92.350	ng	0.00
79) Terphenyl-d14	19.881	244	9463878	88.047	ng	0.00

Target Compounds Qvalue

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

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