

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092023\
 Data File : BM041944.D
 Acq On : 20 Sep 2023 18:39
 Operator : MA/JU
 Sample : 04436-06
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CPT103-05

Quant Time: Sep 21 03:16:17 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM091823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 19 04:49:34 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	291299	20.000	ng	-0.01
21) Naphthalene-d8	10.775	136	1214074	20.000	ng	-0.01
39) Acenaphthene-d10	14.610	164	781802	20.000	ng	0.00
64) Phenanthrene-d10	17.356	188	1667558	20.000	ng	-0.01
76) Chrysene-d12	21.527	240	1375135	20.000	ng	-0.02
86) Perylene-d12	23.944	264	1230019	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	1506251	86.125	ng	0.00
7) Phenol-d6	7.104	99	2070659	85.272	ng	0.00
23) Nitrobenzene-d5	9.151	82	1259638	52.826	ng	-0.01
42) 2,4,6-Tribromophenol	16.098	330	794082	81.406	ng	-0.01
45) 2-Fluorobiphenyl	13.227	172	2669218	50.112	ng	0.00
79) Terphenyl-d14	19.968	244	3688443	48.851	ng	0.00

Target Compounds	Qvalue

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

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