

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092023\
 Data File : BM041948.D
 Acq On : 20 Sep 2023 21:06
 Operator : MA/JU
 Sample : 04453-06
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B102-07

Quant Time: Sep 21 03:17:06 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	295625	20.000	ng	-0.01
21) Naphthalene-d8	10.775	136	1218638	20.000	ng	-0.01
39) Acenaphthene-d10	14.604	164	769263	20.000	ng	-0.01
64) Phenanthrene-d10	17.357	188	1616661	20.000	ng	-0.01
76) Chrysene-d12	21.527	240	1227666	20.000	ng	-0.02
86) Perylene-d12	23.945	264	1109337	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	2128976	119.950	ng	0.00
7) Phenol-d6	7.105	99	2906176	117.928	ng	0.00
23) Nitrobenzene-d5	9.151	82	1752248	73.210	ng	-0.01
42) 2,4,6-Tribromophenol	16.098	330	1049994	109.396	ng	-0.01
45) 2-Fluorobiphenyl	13.222	172	3783492	72.190	ng	-0.01
79) Terphenyl-d14	19.968	244	5095220	75.590	ng	0.00

Target Compounds	Qvalue

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

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