

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091923\
 Data File : BM041929.D
 Acq On : 20 Sep 2023 07:05
 Operator : MA/JU
 Sample : 04453-11
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 FB-9-14

Quant Time: Sep 20 07:39:14 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	262540	20.000	ng	-0.01
21) Naphthalene-d8	10.775	136	1094788	20.000	ng	-0.01
39) Acenaphthene-d10	14.604	164	721167	20.000	ng	-0.01
64) Phenanthrene-d10	17.357	188	1595581	20.000	ng	-0.01
76) Chrysene-d12	21.527	240	1420913	20.000	ng	-0.02
86) Perylene-d12	23.944	264	1321448	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	796721	50.545	ng	0.00
7) Phenol-d6	7.098	99	603748	27.586	ng	-0.01
23) Nitrobenzene-d5	9.151	82	1929049	89.714	ng	-0.01
42) 2,4,6-Tribromophenol	16.098	330	1341459	149.084	ng	-0.01
45) 2-Fluorobiphenyl	13.227	172	4422984	90.019	ng	0.00
79) Terphenyl-d14	19.968	244	7789983	99.850	ng	0.00

Target Compounds	Qvalue

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

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