

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092023\
 Data File : BM041936.D
 Acq On : 20 Sep 2023 13:45
 Operator : MA/JU
 Sample : 04343-02RE
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SAMPLE-2-23RD-STREETRE

Quant Time: Sep 20 14:19:50 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	312016	20.000	ng	-0.01
21) Naphthalene-d8	10.775	136	1304709	20.000	ng	-0.01
39) Acenaphthene-d10	14.604	164	852346	20.000	ng	-0.01
64) Phenanthrene-d10	17.357	188	1852189	20.000	ng	-0.01
76) Chrysene-d12	21.533	240	1645963	20.000	ng	-0.01
86) Perylene-d12	23.944	264	1539654	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	919674	49.094	ng	0.00
7) Phenol-d6	7.099	99	794724	30.555	ng	-0.01
23) Nitrobenzene-d5	9.157	82	2429141	94.795	ng	0.00
42) 2,4,6-Tribromophenol	16.098	330	1044794	98.244	ng	-0.01
45) 2-Fluorobiphenyl	13.228	172	5405126	93.078	ng	0.00
79) Terphenyl-d14	19.968	244	8174743	90.455	ng	0.00

Target Compounds	Qvalue

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

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