

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091923\
 Data File : BM041923.D
 Acq On : 20 Sep 2023 03:29
 Operator : MA/JU
 Sample : 04397-09
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 FB-9-11

Quant Time: Sep 20 04:37:21 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	7.963	152	297099	20.000	ng	0.00
21) Naphthalene-d8	10.781	136	1226984	20.000	ng	0.00
39) Acenaphthene-d10	14.610	164	808046	20.000	ng	0.00
64) Phenanthrene-d10	17.362	188	1692527	20.000	ng	0.00
76) Chrysene-d12	21.527	240	1251859	20.000	ng	-0.02
86) Perylene-d12	23.944	264	1185937	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	973783	54.592	ng	0.00
7) Phenol-d6	7.104	99	791461	31.957	ng	0.00
23) Nitrobenzene-d5	9.157	82	2076367	86.161	ng	0.00
42) 2,4,6-Tribromophenol	16.104	330	1348087	133.712	ng	0.00
45) 2-Fluorobiphenyl	13.227	172	4759934	86.461	ng	0.00
79) Terphenyl-d14	19.968	244	7077711	102.972	ng	0.00

Target Compounds	Qvalue

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

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