

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091923\
 Data File : BM041922.D
 Acq On : 20 Sep 2023 02:53
 Operator : MA/JU
 Sample : 04394-11
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 FB9/8

Quant Time: Sep 20 03:40:25 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	265744	20.000	ng	0.00
21) Naphthalene-d8	10.780	136	1115787	20.000	ng	0.00
39) Acenaphthene-d10	14.610	164	743555	20.000	ng	0.00
64) Phenanthrene-d10	17.356	188	1629888	20.000	ng	-0.01
76) Chrysene-d12	21.533	240	1433838	20.000	ng	-0.01
86) Perylene-d12	23.944	264	1375928	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	1004237	62.943	ng	0.00
7) Phenol-d6	7.104	99	809474	36.541	ng	0.00
23) Nitrobenzene-d5	9.157	82	2012781	91.847	ng	0.00
42) 2,4,6-Tribromophenol	16.104	330	1363612	146.983	ng	0.00
45) 2-Fluorobiphenyl	13.227	172	4546813	89.753	ng	0.00
79) Terphenyl-d14	19.968	244	7590996	96.423	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.410	88	14032	2.156	ng	Qvalue # 87

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

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