

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060923\
 Data File : BM040189.D
 Acq On : 09 Jun 2023 13:39
 Operator : MA/JU
 Sample : 03039-09
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-34

Quant Time: Jun 09 17:10:43 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 08 23:25:50 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.987	152	409513	20.000	ng	0.00
21) Naphthalene-d8	10.804	136	1784039	20.000	ng	0.00
39) Acenaphthene-d10	14.627	164	1064509	20.000	ng	0.00
64) Phenanthrene-d10	17.368	188	2252105	20.000	ng	0.00
76) Chrysene-d12	21.544	240	1758734	20.000	ng	0.00
86) Perylene-d12	23.980	264	1639005	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.540	112	944722	34.446	ng	0.00
7) Phenol-d6	7.140	99	746677	19.952	ng	0.00
23) Nitrobenzene-d5	9.157	82	2299838	69.774	ng	0.00
42) 2,4,6-Tribromophenol	16.110	330	1131395	70.641	ng	0.00
45) 2-Fluorobiphenyl	13.251	172	6045070	75.057	ng	0.00
79) Terphenyl-d14	19.986	244	10039132	105.224	ng	0.00

Target Compounds	Qvalue

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

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