

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061323\
 Data File : BM040231.D
 Acq On : 13 Jun 2023 18:49
 Operator : MA/JU
 Sample : 03047-09
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-41

Quant Time: Jun 13 22:55:53 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM061323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 13 22:36:11 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	569414	20.000	ng	0.00
21) Naphthalene-d8	10.792	136	2372594	20.000	ng	0.00
39) Acenaphthene-d10	14.616	164	1342340	20.000	ng	0.00
64) Phenanthrene-d10	17.362	188	2581203	20.000	ng	0.00
76) Chrysene-d12	21.539	240	1891665	20.000	ng	0.00
86) Perylene-d12	23.968	264	1965108	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.528	112	874718	23.818	ng	0.00
7) Phenol-d6	7.134	99	677902	13.539	ng	0.00
23) Nitrobenzene-d5	9.145	82	2219348	48.981	ng	0.00
42) 2,4,6-Tribromophenol	16.104	330	689413	36.816	ng	0.00
45) 2-Fluorobiphenyl	13.245	172	4790487	48.985	ng	0.00
79) Terphenyl-d14	19.980	244	5978352	61.409	ng	0.00

Target Compounds	Qvalue

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

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