

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM052622\
 Data File : BM035279.D
 Acq On : 26 May 2022 21:53
 Operator : CG/JU
 Sample : N2968-09DL 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-39DL

Quant Time: May 27 03:35:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM052622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 26 18:41:37 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.881	152	107101	20.000	ng	0.00
21) Naphthalene-d8	10.680	136	457341	20.000	ng	# 0.00
39) Acenaphthene-d10	14.515	164	345008	20.000	ng	0.00
64) Phenanthrene-d10	17.262	188	837923	20.000	ng	# 0.00
76) Chrysene-d12	21.444	240	993278	20.000	ng	# 0.00
86) Perylene-d12	23.815	264	1045395	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	56269	10.154	ng	0.00
7) Phenol-d6	7.051	99	50234	6.529	ng	0.00
23) Nitrobenzene-d5	9.039	82	125722	15.847	ng	0.00
42) 2,4,6-Tribromophenol	16.004	330	155539	27.569	ng	0.00
45) 2-Fluorobiphenyl	13.145	172	436212	17.964	ng	0.00
79) Terphenyl-d14	19.886	244	887741	17.503	ng	0.00
Target Compounds						
						Qvalue
27) 2,4-Dimethylphenol	9.863	122	27763	4.907	ng	98
31) Naphthalene	10.733	128	1040162	46.256	ng	100
37) 2-Methylnaphthalene	12.345	142	394405	23.361	ng	98
38) 1-Methylnaphthalene	12.563	142	294244	17.695	ng	100

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

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