

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110923\
 Data File : BM042673.D
 Acq On : 09 Nov 2023 16:06
 Operator : MA/JU
 Sample : 05256-12
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-10

Quant Time: Nov 09 16:40:09 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM103023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 09 13:21:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.845	152	29337	20.000	ng	0.00
21) Naphthalene-d8	10.663	136	100896	20.000	ng	# 0.00
39) Acenaphthene-d10	14.504	164	60852	20.000	ng	0.00
64) Phenanthrene-d10	17.257	188	123064	20.000	ng	# 0.00
76) Chrysene-d12	21.439	240	117736	20.000	ng	# 0.00
86) Perylene-d12	23.815	264	143642	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.381	112	269518	148.267	ng	0.00
7) Phenol-d6	7.010	99	316615	127.000	ng	0.00
23) Nitrobenzene-d5	9.045	82	253546	117.144	ng	0.00
42) 2,4,6-Tribromophenol	15.998	330	139197	175.206	ng	0.00
45) 2-Fluorobiphenyl	13.121	172	490077	108.369	ng	0.00
79) Terphenyl-d14	19.874	244	781449	112.542	ng	0.00

Target Compounds	Qvalue

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

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