

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101023\
 Data File : BM042251.D
 Acq On : 10 Oct 2023 14:03
 Operator : MA/JU
 Sample : PB156171BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK171

Quant Time: Oct 10 15:48:15 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM100223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 10 14:01:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.026	152	4546	0.400	ng/ul	0.00
4) Naphthalene-d8	10.845	136	11876	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.666	164	4770	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	17.417	188	9237	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.600	240	3776	0.400	ng/ul	# 0.00
23) Perylene-d12	24.076	264	3672	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.376	96	40888	6.754	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.429	152	4630	0.284	ng/ul	0.00
18) Fluoranthene-d10	19.441	212	5255	0.371	ng/ul	0.00
Target Compounds						
					Qvalue	
12) Fluorene	15.661	166	1986	0.077	ng/ul#	15
20) Pyrene	19.808	202	1617	0.053	ng/ul#	1
26) Benzo(a)pyrene	23.915	252	2783	0.149	ng/ul#	56

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

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