

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100923\
 Data File : BM042230.D
 Acq On : 09 Oct 2023 18:57
 Operator : MA/JU
 Sample : PB156171BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK171

Quant Time: Oct 10 00:01:13 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 : Mon Oct 09 23:59:28 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.043	152	5	0.400	ng/ul	# 0.00
4) Naphthalene-d8	10.878	136	90	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.694	164	240	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	17.447	188	360	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.638	240	180	0.400	ng/ul	# 0.00
23) Perylene-d12	0.000	264	0	0.000	ng/ul	-24.14
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.410	96	40806	6128.451	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.456	152	4523	36.642	ng/ul	0.00
18) Fluoranthene-d10	19.469	212	5883	8.705	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.928	128	88	0.293	ng/ul#	1
7) 2-Methylnaphthalene	12.506	142	4	0.022	ng/ul	87
8) 1-Methylnaphthalene	12.742	142	35	0.193	ng/ul#	76
12) Fluorene	15.684	166	1989	1.538	ng/ul#	16
14) Pentachlorophenol	17.084	266	18	0.110	ng/ul#	55
15) Phenanthrene	17.485	178	48	0.032	ng/ul#	1
16) Anthracene	17.544	178	438	0.323	ng/ul#	1
19) Fluoranthene	19.501	202	50	0.037	ng/ul#	1
20) Pyrene	19.840	202	1630	1.122	ng/ul#	1
21) Benzo(a)anthracene	21.621	228	97	0.083	ng/ul#	2
22) Chrysene	21.679	228	105	0.093	ng/ul#	18

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

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