

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090923\
 Data File : BM041758.D
 Acq On : 09 Sep 2023 20:34
 Operator : MA/JU
 Sample : 04247-09
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BBKB6

Quant Time: Sep 09 23:34:19 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM090523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 09 23:26:58 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.001	152	131	0.400	ng/ul	# 0.00
4) Naphthalene-d8	10.812	136	294	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.629	164	113	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	17.379	188	212	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.548	240	75	0.400	ng/ul	# 0.00
23) Perylene-d12	23.971	264	51	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.427	96	577	2.882	ng/ul	0.03
6) 2-Methylnaphthalene-d10	12.385	152	143	0.378	ng/ul	0.00
18) Fluoranthene-d10	19.399	212	123	0.338	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.473	88	11	0.053	ng/ul#	24
5) Naphthalene	10.856	128	47	0.049	ng/ul#	1
8) 1-Methylnaphthalene	12.682	142	12	0.022	ng/ul#	75
11) Acenaphthene	14.689	153	81	0.149	ng/ul#	76
14) Pentachlorophenol	16.991	266	3	0.035	ng/ul#	1
15) Phenanthrene	17.426	178	19	0.021	ng/ul#	1
19) Fluoranthene	19.431	202	23	0.026	ng/ul#	1
20) Pyrene	19.766	202	59	0.067	ng/ul#	1
22) Chrysene	21.583	228	16	0.024	ng/ul#	1
24) Benzo(b)fluoranthene	23.217	252	19	0.042	ng/ul#	1
25) Benzo(k)fluoranthene	23.278	252	23	0.051	ng/ul#	1
26) Benzo(a)pyrene	23.854	252	42	0.113	ng/ul#	1

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

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