

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100923\
 Data File : BM042246.D
 Acq On : 10 Oct 2023 04:36
 Operator : MA/JU
 Sample : PB156171BS
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS171

Quant Time: Oct 10 05:07:16 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	8.047	152	4	0.400	ng/ul	# 0.00
4) Naphthalene-d8	10.873	136	97	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.689	164	304	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	17.443	188	420	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.626	240	177	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	37926	7119.897	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.451	152	5378	40.425	ng/ul	0.00
18) Fluoranthene-d10	19.464	212	7554	11.367	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.444	88	80108	14401.031	ng/ul#	73
5) Naphthalene	10.928	128	1260841	3896.207	ng/ul	98
7) 2-Methylnaphthalene	12.528	142	787020	4061.868	ng/ul	99
8) 1-Methylnaphthalene	12.748	142	751187	3848.770	ng/ul	99
10) Acenaphthylene	14.421	152	1328499	691.794	ng/ul	99
11) Acenaphthene	14.754	153	855488	588.512	ng/ul	98
12) Fluorene	15.739	166	950084	580.157	ng/ul	99
14) Pentachlorophenol	17.071	266	292982	1538.759	ng/ul	98
15) Phenanthrene	17.485	178	1462382	844.521	ng/ul	100
16) Anthracene	17.578	178	1429202	902.276	ng/ul	100
19) Fluoranthene	19.496	202	1619818	1204.743	ng/ul	100
20) Pyrene	19.864	202	1559868	1091.531	ng/ul	98
21) Benzo(a)anthracene	21.609	228	1196690	1047.286	ng/ul	98
22) Chrysene	21.664	228	1082075	972.938	ng/ul	99

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

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