

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071723\
 Data File : BM040946.D
 Acq On : 18 Jul 2023 09:05
 Operator : MA/JU
 Sample : PB153976BS
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS976

Quant Time: Jul 18 22:25:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 18 22:24:09 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.853	152	4530	0.400	ng/ul	0.00
4) Naphthalene-d8	10.653	136	15182	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.499	164	6268	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.253	188	10301	0.400	ng/ul	0.00
17) Chrysene-d12	21.448	240	6740	0.400	ng/ul	0.00
23) Perylene-d12	23.822	264	7719	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	4690	0.659	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.247	152	6614	0.311	ng/ul	0.00
18) Fluoranthene-d10	19.283	212	7425	0.344	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.309	88	10548	1.456	ng/ul	90
5) Naphthalene	10.702	128	11737	0.280	ng/ul	98
7) 2-Methylnaphthalene	12.319	142	6383	0.260	ng/ul	97
8) 1-Methylnaphthalene	12.539	142	6780	0.276	ng/ul	98
10) Acenaphthylene	14.222	152	7459	0.266	ng/ul	98
11) Acenaphthene	14.560	153	6373	0.270	ng/ul	98
12) Fluorene	15.554	166	6285	0.249	ng/ul	100
14) Pentachlorophenol	16.911	266	1037	0.523	ng/ul	94
15) Phenanthrene	17.295	178	8389	0.262	ng/ul	98
16) Anthracene	17.388	178	6757	0.251	ng/ul	98
19) Fluoranthene	19.315	202	8170	0.285	ng/ul	100
20) Pyrene	19.678	202	8490	0.277	ng/ul	97
21) Benzo(a)anthracene	21.431	228	6525	0.286	ng/ul	99
22) Chrysene	21.483	228	7217	0.278	ng/ul	99
24) Benzo(b)fluoranthene	23.097	252	7850	0.263	ng/ul	93
25) Benzo(k)fluoranthene	23.149	252	7794	0.265	ng/ul#	92
26) Benzo(a)pyrene	23.716	252	7236	0.269	ng/ul	90
27) Indeno(1,2,3-cd)pyrene	26.280	276	10499	0.280	ng/ul	96
28) Dibenzo(a,h)anthracene	26.294	278	8348	0.285	ng/ul	96
29) Benzo(g,h,i)perylene	27.031	276	9196	0.280	ng/ul	99

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

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