

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072123\
 Data File : BM041015.D
 Acq On : 21 Jul 2023 13:07
 Operator : MA/JU
 Sample : PB154070BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS070

Quant Time: Jul 21 16:40:56 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM071923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 21 16:33:57 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.870	152	3238	0.400	ng/ul	0.00
4) Naphthalene-d8	10.675	136	12168	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.518	164	7110	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.278	188	14241	0.400	ng/ul	0.00
17) Chrysene-d12	21.469	240	11228	0.400	ng/ul	0.00
23) Perylene-d12	23.863	264	9788	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.280	96	3914	0.735	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.269	152	7271	0.448	ng/ul	0.00
18) Fluoranthene-d10	19.306	212	16064	0.340	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.313	88	9151	1.633	ng/ul#	71
5) Naphthalene	10.724	128	11941	0.359	ng/ul	100
7) 2-Methylnaphthalene	12.346	142	7088	0.388	ng/ul	100
8) 1-Methylnaphthalene	12.561	142	7912	0.408	ng/ul	100
10) Acenaphthylene	14.240	152	10665	0.358	ng/ul	99
11) Acenaphthene	14.583	153	9014	0.343	ng/ul	97
12) Fluorene	15.577	166	9812	0.361	ng/ul	98
14) Pentachlorophenol	16.936	266	2347	0.818	ng/ul	99
15) Phenanthrene	17.316	178	15145	0.356	ng/ul	99
16) Anthracene	17.413	178	12168	0.367	ng/ul	99
19) Fluoranthene	19.338	202	17881	0.308	ng/ul	94
20) Pyrene	19.701	202	19104	0.323	ng/ul	94
21) Benzo(a)anthracene	21.454	228	13118	0.384	ng/ul	100
22) Chrysene	21.504	228	14819	0.379	ng/ul	100
24) Benzo(b)fluoranthene	23.132	252	13852	0.369	ng/ul	95
25) Benzo(k)fluoranthene	23.178	252	13314	0.373	ng/ul#	95
26) Benzo(a)pyrene	23.757	252	11854	0.365	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	26.337	276	15843	0.368	ng/ul#	90
28) Dibenzo(a,h)anthracene	26.354	278	12386	0.380	ng/ul	94
29) Benzo(g,h,i)perylene	27.095	276	14143	0.355	ng/ul	94

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

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