

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071423\
 Data File : BM040892.D
 Acq On : 14 Jul 2023 23:19
 Operator : MA/JU
 Sample : 03437-11MS
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A46K7MS

Quant Time: Jul 15 04:00:35 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 : Sat Jul 15 03:58:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.861	152	4569	0.400	ng/ul	0.00
4) Naphthalene-d8	10.664	136	14523	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.509	164	6045	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.261	188	10913	0.400	ng/ul	0.00
17) Chrysene-d12	21.448	240	7436	0.400	ng/ul	0.00
23) Perylene-d12	23.825	264	7350	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.284	96	4065	0.566	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.258	152	6238	0.306	ng/ul	0.00
18) Fluoranthene-d10	19.292	212	8657	0.363	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.317	88	9199	1.259	ng/ul	90
5) Naphthalene	10.713	128	12021	0.300	ng/ul	98
7) 2-Methylnaphthalene	12.330	142	6755	0.287	ng/ul	100
8) 1-Methylnaphthalene	12.550	142	7045	0.300	ng/ul	98
10) Acenaphthylene	14.226	152	8111	0.300	ng/ul	98
11) Acenaphthene	14.569	153	6719	0.295	ng/ul	99
12) Fluorene	15.559	166	6896	0.283	ng/ul	99
14) Pentachlorophenol	16.919	266	438	0.208	ng/ul	93
15) Phenanthrene	17.303	178	12424	0.367	ng/ul	99
16) Anthracene	17.392	178	8682	0.305	ng/ul	97
19) Fluoranthene	19.320	202	14372	0.455	ng/ul	99
20) Pyrene	19.682	202	14906	0.440	ng/ul	99
21) Benzo(a)anthracene	21.434	228	10025	0.399	ng/ul	99
22) Chrysene	21.483	228	10708	0.373	ng/ul	98
24) Benzo(b)fluoranthene	23.099	252	11436	0.402	ng/ul	98
25) Benzo(k)fluoranthene	23.149	252	9787	0.349	ng/ul	95
26) Benzo(a)pyrene	23.719	252	9357	0.365	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.280	276	12108	0.339	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.300	278	9005	0.322	ng/ul	97
29) Benzo(g,h,i)perylene	27.038	276	10659	0.341	ng/ul	98

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

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