

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070823\
 Data File : BM040773.D
 Acq On : 09 Jul 2023 12:38
 Operator : MA/JU
 Sample : 03356-10MS
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 YBW38MS

Quant Time: Jul 09 14:44:09 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 : Sun Jul 09 14:07:58 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.891	152	4226	0.400	ng/ul	0.00
4) Naphthalene-d8	10.691	136	15409	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.536	164	6780	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.286	188	11218	0.400	ng/ul	0.00
17) Chrysene-d12	21.477	240	6729	0.400	ng/ul	0.00
23) Perylene-d12	23.874	264	5864	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	4010	0.604	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.286	152	6051	0.280	ng/ul	0.00
18) Fluoranthene-d10	19.315	212	7982	0.370	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.334	88	9995	1.479	ng/ul	91
5) Naphthalene	10.741	128	12343	0.290	ng/ul	98
7) 2-Methylnaphthalene	12.357	142	7212	0.289	ng/ul	97
8) 1-Methylnaphthalene	12.577	142	6812	0.273	ng/ul	98
10) Acenaphthylene	14.254	152	8919	0.294	ng/ul	95
11) Acenaphthene	14.597	153	7421	0.290	ng/ul	98
12) Fluorene	15.587	166	7843	0.287	ng/ul	98
14) Pentachlorophenol	16.949	266	406	0.188	ng/ul	90
15) Phenanthrene	17.329	178	11049	0.317	ng/ul	98
16) Anthracene	17.421	178	10574	0.361	ng/ul	94
19) Fluoranthene	19.348	202	10329	0.361	ng/ul	98
20) Pyrene	19.710	202	10943	0.357	ng/ul	97
21) Benzo(a)anthracene	21.463	228	7077	0.311	ng/ul	97
22) Chrysene	21.515	228	8011	0.309	ng/ul	99
24) Benzo(b)fluoranthene	23.143	252	7014	0.309	ng/ul	84
25) Benzo(k)fluoranthene	23.190	252	7677	0.343	ng/ul#	85
26) Benzo(a)pyrene	23.766	252	6333	0.310	ng/ul#	77
27) Indeno(1,2,3-cd)pyrene	26.354	276	7946	0.279	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.374	278	6261	0.281	ng/ul#	87
29) Benzo(g,h,i)perylene	27.122	276	7371	0.295	ng/ul	96

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

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