

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070523\
 Data File : BM040727.D
 Acq On : 07 Jul 2023 23:59
 Operator : MA/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4161

Quant Time: Jul 08 05:11:40 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 : Fri Jul 07 10:22:24 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	4822	0.400	ng/ul	0.00
4) Naphthalene-d8	10.708	136	18271	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.546	164	8541	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.299	188	15081	0.400	ng/ul	0.00
17) Chrysene-d12	21.486	240	9342	0.400	ng/ul	0.00
23) Perylene-d12	23.886	264	8148	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	2950	0.389	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.297	152	9684	0.378	ng/ul	0.00
18) Fluoranthene-d10	19.329	212	12720	0.425	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.343	88	2621	0.340	ng/ul#	92
5) Naphthalene	10.757	128	18620	0.369	ng/ul	99
7) 2-Methylnaphthalene	12.368	142	10969	0.371	ng/ul	97
8) 1-Methylnaphthalene	12.588	142	10344	0.350	ng/ul	98
10) Acenaphthylene	14.264	152	13610	0.357	ng/ul	98
11) Acenaphthene	14.606	153	11618	0.361	ng/ul	98
12) Fluorene	15.601	166	12322	0.358	ng/ul	99
14) Pentachlorophenol	16.957	266	415	0.143	ng/ul#	86
15) Phenanthrene	17.341	178	16469	0.352	ng/ul	98
16) Anthracene	17.438	178	14309	0.363	ng/ul	97
19) Fluoranthene	19.357	202	15908	0.401	ng/ul	99
20) Pyrene	19.719	202	16392	0.385	ng/ul	98
21) Benzo(a)anthracene	21.469	228	11004	0.348	ng/ul	100
22) Chrysene	21.521	228	12263	0.340	ng/ul	99
24) Benzo(b)fluoranthene	23.152	252	11791	0.374	ng/ul	94
25) Benzo(k)fluoranthene	23.196	252	11831	0.381	ng/ul	94
26) Benzo(a)pyrene	23.778	252	10228	0.360	ng/ul	91
27) Indeno(1,2,3-cd)pyrene	26.374	276	13151	0.332	ng/ul	95
28) Dibenzo(a,h)anthracene	26.387	278	10339	0.334	ng/ul	95
29) Benzo(g,h,i)perylene	27.135	276	11814	0.341	ng/ul	99

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

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