

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090623\
 Data File : BM041699.D
 Acq On : 07 Sep 2023 17:41
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4072

Quant Time: Sep 08 02:54:10 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM090523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 06 02:20:07 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.000	152	5751	0.400	ng/ul	0.00
4) Naphthalene-d8	10.817	136	13474	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.643	164	5388	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.392	188	10207	0.400	ng/ul	0.00
17) Chrysene-d12	21.562	240	4764	0.400	ng/ul	0.00
23) Perylene-d12	23.994	264	5659	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.406	96	3062	0.348	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.396	152	6775	0.391	ng/ul	-0.01
18) Fluoranthene-d10	19.413	212	7949	0.344	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.439	88	3243	0.355	ng/ul#	86
5) Naphthalene	10.867	128	17075	0.386	ng/ul	100
7) 2-Methylnaphthalene	12.473	142	9929	0.395	ng/ul	99
8) 1-Methylnaphthalene	12.693	142	9950	0.394	ng/ul	99
10) Acenaphthylene	14.370	152	13217	0.397	ng/ul	99
11) Acenaphthene	14.703	153	9981	0.386	ng/ul	99
12) Fluorene	15.693	166	10898	0.375	ng/ul	99
14) Pentachlorophenol	17.016	266	1719	0.421	ng/ul	98
15) Phenanthrene	17.438	178	16453	0.382	ng/ul	99
16) Anthracene	17.527	178	14799	0.401	ng/ul	99
19) Fluoranthene	19.445	202	17573	0.313	ng/ul	99
20) Pyrene	19.808	202	18422	0.330	ng/ul	99
21) Benzo(a)anthracene	21.544	228	14476	0.373	ng/ul	99
22) Chrysene	21.600	228	14316	0.342	ng/ul	99
24) Benzo(b)fluoranthene	23.243	252	15179	0.300	ng/ul	95
25) Benzo(k)fluoranthene	23.295	252	15106	0.305	ng/ul	96
26) Benzo(a)pyrene	23.886	252	13343	0.324	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.515	276	18289	0.324	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.535	278	12908	0.326	ng/ul	99
29) Benzo(g,h,i)perylene	27.300	276	15631	0.313	ng/ul	99

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

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