

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062623\
 Data File : BM040396.D
 Acq On : 26 Jun 2023 11:42
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4132

Quant Time: Jun 27 04:25:02 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 Jun 23 00:44:24 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.949	152	5934	0.400	ng/ul	0.00
4) Naphthalene-d8	10.756	136	19203	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.586	164	8100	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.332	188	13567	0.400	ng/ul	0.00
17) Chrysene-d12	21.515	240	9498	0.400	ng/ul	0.00
23) Perylene-d12	23.929	264	10029	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.337	96	3666	0.393	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.345	152	10379	0.386	ng/ul	0.00
18) Fluoranthene-d10	19.361	212	11867	0.390	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.371	88	3856	0.406	ng/ul	87
5) Naphthalene	10.805	128	21530	0.406	ng/ul	100
7) 2-Methylnaphthalene	12.417	142	11925	0.383	ng/ul	99
8) 1-Methylnaphthalene	12.637	142	11819	0.380	ng/ul	99
10) Acenaphthylene	14.309	152	13929	0.385	ng/ul	99
11) Acenaphthene	14.651	153	11878	0.389	ng/ul	98
12) Fluorene	15.636	166	11934	0.366	ng/ul	99
14) Pentachlorophenol	16.990	266	629	0.241	ng/ul	97
15) Phenanthrene	17.374	178	16130	0.383	ng/ul	99
16) Anthracene	17.467	178	13438	0.379	ng/ul	98
19) Fluoranthene	19.389	202	15452	0.383	ng/ul	100
20) Pyrene	19.751	202	16322	0.377	ng/ul	98
21) Benzo(a)anthracene	21.497	228	12511	0.389	ng/ul	99
22) Chrysene	21.550	228	14002	0.382	ng/ul	100
24) Benzo(b)fluoranthene	23.189	252	15176	0.391	ng/ul	93
25) Benzo(k)fluoranthene	23.239	252	14655	0.383	ng/ul	96
26) Benzo(a)pyrene	23.821	252	13844	0.396	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	26.437	276	18924	0.389	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.454	278	14756	0.387	ng/ul	99
29) Benzo(g,h,i)perylene	27.205	276	16159	0.379	ng/ul	100

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

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