

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060321\
 Data File : BM030282.D
 Acq On : 03 Jun 2021 13:46
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4088

Quant Time: Jun 03 14:21:29 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM060121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 03 10:49:01 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.860	152	5410	0.40	ng/ul	0.00
4) Naphthalene-d8	10.648	136	18536	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.481	164	9978	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.214	188	19247	0.40	ng/ul	0.00
17) Chrysene-d12	21.376	240	12377	0.40	ng/ul	0.00
23) Perylene-d12	23.663	264	9534	0.40	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	2249	0.42	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.236	152	11120	0.42	ng/ul	0.00
18) Fluoranthene-d10	19.233	212	18354	0.44	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.365	88	2296	0.40	ng/ul	94
5) Naphthalene	10.698	128	22292	0.41	ng/ul	100
7) 2-Methylnaphthalene	12.308	142	15117	0.41	ng/ul	99
8) 1-Methylnaphthalene	12.528	142	14494	0.41	ng/ul	99
10) Acenaphthylene	14.201	152	17569	0.37	ng/ul	100
11) Acenaphthene	14.542	153	15481	0.38	ng/ul	98
12) Fluorene	15.528	166	17491	0.38	ng/ul	100
14) Pentachlorophenol	16.867	266	1942	0.32	ng/ul	99
15) Phenanthrene	17.256	178	27078	0.39	ng/ul	100
16) Anthracene	17.349	178	22666	0.38	ng/ul	99
19) Fluoranthene	19.264	202	27193	0.43	ng/ul	100
20) Pyrene	19.626	202	26511	0.41	ng/ul	98
21) Benzo(a)anthracene	21.360	228	18700	0.38	ng/ul	100
22) Chrysene	21.410	228	21236	0.40	ng/ul	100
24) Benzo(b)fluoranthene	22.970	252	20014	0.42	ng/ul	97
25) Benzo(k)fluoranthene	23.016	252	18918	0.42	ng/ul	97
26) Benzo(a)pyrene	23.561	252	15182	0.40	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.986	276	19656	0.39	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.996	278	15850	0.39	ng/ul	99
29) Benzo(g,h,i)perylene	26.696	276	17364	0.39	ng/ul	99

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

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