

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062122\
 Data File : BN020316.D
 Acq On : 21 Jun 2022 20:52
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4375

Quant Time: Jun 22 01:22:11 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN061822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 20 01:13:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.800	152	581	0.400	ng/ul	0.00
4) Naphthalene-d8	10.584	136	2060	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.423	164	1505	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.166	188	3929	0.400	ng/ul	0.00
17) Chrysene-d12	21.349	240	6762	0.400	ng/ul	0.00
23) Perylene-d12	23.655	264	8291	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.256	96	265	0.399	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.179	152	1248	0.366	ng/ul	0.00
18) Fluoranthene-d10	19.192	212	5722	0.410	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.294	88	220	0.345	ng/ul	90
5) Naphthalene	10.634	128	2302	0.349	ng/ul	97
7) 2-Methylnaphthalene	12.256	142	1381	0.324	ng/ul	99
8) 1-Methylnaphthalene	12.470	142	1626	0.388	ng/ul	98
10) Acenaphthylene	14.141	152	2567	0.357	ng/ul	97
11) Acenaphthene	14.483	153	2048	0.298	ng/ul	99
12) Fluorene	15.473	166	2497	0.350	ng/ul	98
14) Pentachlorophenol	16.837	266	288	0.382	ng/ul	99
15) Phenanthrene	17.208	178	4898	0.344	ng/ul	98
16) Anthracene	17.301	178	4081	0.339	ng/ul	98
19) Fluoranthene	19.225	202	7042	0.258	ng/ul	99
20) Pyrene	19.583	202	7801	0.264	ng/ul	99
21) Benzo(a)anthracene	21.334	228	7873	0.300	ng/ul	99
22) Chrysene	21.384	228	9884	0.302	ng/ul	99
24) Benzo(b)fluoranthene	22.959	252	12311	0.208	ng/ul	98
25) Benzo(k)fluoranthene	23.003	252	12586	0.228	ng/ul	99
26) Benzo(a)pyrene	23.555	252	12485	0.333	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.029	276	17089	0.280	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.042	278	13707	0.282	ng/ul	99
29) Benzo(g,h,i)perylene	26.750	276	15673	0.291	ng/ul	100

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

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