

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032223\
 Data File : BN024504.D
 Acq On : 23 Mar 2023 07:19
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4252

Quant Time: Mar 23 07:48:38 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN030923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 23 03:32:22 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.025	152	12566	0.400	ng/ul	0.00
4) Naphthalene-d8	10.839	136	37184	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.661	164	20465	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.405	188	42738	0.400	ng/ul	0.00
17) Chrysene-d12	21.597	240	36230	0.400	ng/ul	0.00
23) Perylene-d12	24.088	264	31746	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	4539	0.328	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.423	152	17979	0.407	ng/ul	0.00
18) Fluoranthene-d10	19.432	212	37657	0.391	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.431	88	4959	0.330	ng/ul#	88
5) Naphthalene	10.889	128	35821	0.386	ng/ul	100
7) 2-Methylnaphthalene	12.495	142	22318	0.396	ng/ul	100
8) 1-Methylnaphthalene	12.715	142	23399	0.395	ng/ul	100
10) Acenaphthylene	14.384	152	32859	0.385	ng/ul	100
11) Acenaphthene	14.721	153	24683	0.375	ng/ul	97
12) Fluorene	15.707	166	27478	0.374	ng/ul	99
14) Pentachlorophenol	17.054	266	3771	0.318	ng/ul#	72
15) Phenanthrene	17.447	178	45228	0.369	ng/ul	100
16) Anthracene	17.540	178	41140	0.383	ng/ul	99
19) Fluoranthene	19.459	202	52357	0.363	ng/ul	100
20) Pyrene	19.822	202	54781	0.374	ng/ul	99
21) Benzo(a)anthracene	21.580	228	46411	0.392	ng/ul	100
22) Chrysene	21.635	228	45831	0.367	ng/ul	99
24) Benzo(b)fluoranthene	23.325	252	44992	0.371	ng/ul	95
25) Benzo(k)fluoranthene	23.372	252	43906	0.349	ng/ul	94
26) Benzo(a)pyrene	23.974	252	38109	0.365	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	26.680	276	42830	0.365	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.700	278	33150	0.371	ng/ul	98
29) Benzo(g,h,i)perylene	27.485	276	36020	0.353	ng/ul	97

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

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