

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102021\
 Data File : BN017012.D
 Acq On : 20 Oct 2021 09:25
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4272

Quant Time: Oct 20 10:48:47 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN101921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 19 14:22:25 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.548	152	2259	0.40	ng/ul	0.00
4) Naphthalene-d8	10.317	136	8634	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.199	164	5068	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.953	188	11345	0.40	ng/ul	0.00
17) Chrysene-d12	21.162	240	11908	0.40	ng/ul	0.00
23) Perylene-d12	23.375	264	11993	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.076	96	991	0.40	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.923	152	4594	0.37	ng/ul	0.00
18) Fluoranthene-d10	18.995	212	11453	0.39	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.109	88	888	0.30	ng/ul#	49
5) Naphthalene	10.366	128	9276	0.35	ng/ul	99
7) 2-Methylnaphthalene	11.994	142	6146	0.38	ng/ul	100
8) 1-Methylnaphthalene	12.220	142	6300	0.38	ng/ul	100
10) Acenaphthylene	13.916	152	7646	0.37	ng/ul	99
11) Acenaphthene	14.259	153	6917	0.39	ng/ul	99
12) Fluorene	15.253	166	7772	0.38	ng/ul	100
14) Pentachlorophenol	16.611	266	715	0.26	ng/ul	99
15) Phenanthrene	16.995	178	14034	0.39	ng/ul	99
16) Anthracene	17.088	178	11138	0.37	ng/ul	100
19) Fluoranthene	19.023	202	16087	0.39	ng/ul	99
20) Pyrene	19.385	202	16407	0.39	ng/ul	99
21) Benzo(a)anthracene	21.148	228	13571	0.35	ng/ul	99
22) Chrysene	21.200	228	17756	0.40	ng/ul	100
24) Benzo(b)fluoranthene	22.711	252	18096	0.40	ng/ul	99
25) Benzo(k)fluoranthene	22.755	252	19965	0.40	ng/ul	99
26) Benzo(a)pyrene	23.278	252	14992	0.37	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.604	276	22490	0.39	ng/ul	100
28) Dibenzo(a,h)anthracene	25.624	278	17807	0.39	ng/ul	100
29) Benzo(g,h,i)perylene	26.278	276	19908	0.40	ng/ul	99

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

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