

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092122\
 Data File : BN021899.D
 Acq On : 23 Sep 2022 13:30
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4312

Quant Time: Sep 23 16:18:43 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN090822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 20 00:20:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.042	152	4921	0.400	ng/u1	0.00
4) Naphthalene-d8	10.872	136	16365	0.400	ng/u1	#-0.01
9) Acenaphthene-d10	14.703	164	9132	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.456	188	16982	0.400	ng/u1	0.00
17) Chrysene-d12	21.650	240	10288	0.400	ng/u1	# 0.00
23) Perylene-d12	24.158	264	8436	0.400	ng/u1	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.359	96	2221	0.367	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.462	152	11060	0.427	ng/u1	-0.01
18) Fluoranthene-d10	19.483	212	17431	0.455	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.397	88	2338	0.377	ng/u1	94
5) Naphthalene	10.922	128	19435	0.421	ng/u1	98
7) 2-Methylnaphthalene	12.533	142	12714	0.424	ng/u1	98
8) 1-Methylnaphthalene	12.753	142	12974	0.422	ng/u1	99
10) Acenaphthylene	14.425	152	16521	0.431	ng/u1	98
11) Acenaphthene	14.768	153	13768	0.424	ng/u1	99
12) Fluorene	15.753	166	15017	0.412	ng/u1	99
14) Pentachlorophenol	17.109	266	815	0.248	ng/u1	100
15) Phenanthrene	17.498	178	22412	0.418	ng/u1	99
16) Anthracene	17.591	178	19240	0.433	ng/u1	98
19) Fluoranthene	19.515	202	21110	0.440	ng/u1	98
20) Pyrene	19.878	202	21284	0.449	ng/u1	96
21) Benzo(a)anthracene	21.633	228	15723	0.456	ng/u1	99
22) Chrysene	21.688	228	15754	0.419	ng/u1	99
24) Benzo(b)fluoranthene	23.384	252	14815	0.362	ng/u1	93
25) Benzo(k)fluoranthene	23.436	252	14935	0.376	ng/u1#	93
26) Benzo(a)pyrene	24.044	252	12510	0.404	ng/u1#	92
27) Indeno(1,2,3-cd)pyrene	26.804	276	14884	0.427	ng/u1#	98
28) Dibenzo(a,h)anthracene	26.821	278	12704	0.429	ng/u1	95
29) Benzo(g,h,i)perylene	27.619	276	13701	0.426	ng/u1	97

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

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