

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092122\
 Data File : BN021859.D
 Acq On : 22 Sep 2022 03:26
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4308

Quant Time: Sep 22 03:58:20 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	8.051	152	5439	0.400 ng/u1	0.00
4) Naphthalene-d8	10.878	136	18091	0.400 ng/u1	# 0.00
9) Acenaphthene-d10	14.707	164	9729	0.400 ng/u1	0.00
13) Phenanthrene-d10	17.455	188	17351	0.400 ng/u1	# 0.00
17) Chrysene-d12	21.647	240	10079	0.400 ng/u1	# 0.00
23) Perylene-d12	24.155	264	8090	0.400 ng/u1	# 0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.363	96	2584	0.386 ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.467	152	11639	0.406 ng/u1	0.00
18) Fluoranthene-d10	19.487	212	16395	0.437 ng/u1	0.00
Target Compounds					
					Qvalue
2) 1,4-Dioxane	3.401	88	2571	0.375 ng/u1	94
5) Naphthalene	10.927	128	20707	0.406 ng/u1	97
7) 2-Methylnaphthalene	12.539	142	13329	0.402 ng/u1	98
8) 1-Methylnaphthalene	12.758	142	13677	0.402 ng/u1	99
10) Acenaphthylene	14.430	152	16316	0.399 ng/u1	98
11) Acenaphthene	14.768	153	14134	0.408 ng/u1	99
12) Fluorene	15.753	166	15208	0.392 ng/u1	98
14) Pentachlorophenol	17.109	266	978	0.291 ng/u1	99
15) Phenanthrene	17.498	178	22138	0.404 ng/u1	99
16) Anthracene	17.591	178	18163	0.400 ng/u1	99
19) Fluoranthene	19.515	202	20035	0.426 ng/u1	97
20) Pyrene	19.878	202	19926	0.429 ng/u1	94
21) Benzo(a)anthracene	21.632	228	14556	0.431 ng/u1	100
22) Chrysene	21.685	228	15264	0.414 ng/u1	100
24) Benzo(b)fluoranthene	23.380	252	14360	0.366 ng/u1	93
25) Benzo(k)fluoranthene	23.433	252	14289	0.376 ng/u1#	94
26) Benzo(a)pyrene	24.044	252	11933	0.401 ng/u1#	91
27) Indeno(1,2,3-cd)pyrene	26.801	276	14260	0.426 ng/u1	97
28) Dibenzo(a,h)anthracene	26.818	278	12092	0.426 ng/u1	95
29) Benzo(g,h,i)perylene	27.609	276	13171	0.427 ng/u1	97

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

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