

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092122\
 Data File : BN021862.D
 Acq On : 22 Sep 2022 06:36
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4309

Quant Time: Sep 22 07:23:48 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	5290	0.400	ng/u1	0.00
4) Naphthalene-d8	10.878	136	17486	0.400	ng/u1	# 0.00
9) Acenaphthene-d10	14.708	164	9296	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.460	188	16812	0.400	ng/u1	0.00
17) Chrysene-d12	21.650	240	9703	0.400	ng/u1	# 0.00
23) Perylene-d12	24.161	264	7426	0.400	ng/u1	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.363	96	2441	0.375	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.467	152	11128	0.402	ng/u1	0.00
18) Fluoranthene-d10	19.487	212	15797	0.438	ng/u1	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.401	88	2475	0.371	ng/u1	94
5) Naphthalene	10.927	128	19994	0.406	ng/u1	97
7) 2-Methylnaphthalene	12.539	142	12796	0.399	ng/u1	98
8) 1-Methylnaphthalene	12.759	142	13090	0.398	ng/u1	99
10) Acenaphthylene	14.425	152	15399	0.394	ng/u1	98
11) Acenaphthene	14.768	153	13538	0.409	ng/u1	99
12) Fluorene	15.758	166	14635	0.394	ng/u1	98
14) Pentachlorophenol	17.109	266	883	0.271	ng/u1	99
15) Phenanthrene	17.502	178	21496	0.405	ng/u1	99
16) Anthracene	17.591	178	17401	0.395	ng/u1	99
19) Fluoranthene	19.515	202	19511	0.431	ng/u1	97
20) Pyrene	19.882	202	19264	0.431	ng/u1	97
21) Benzo(a)anthracene	21.635	228	13732	0.422	ng/u1	100
22) Chrysene	21.688	228	14666	0.414	ng/u1	100
24) Benzo(b)fluoranthene	23.383	252	13746	0.381	ng/u1	94
25) Benzo(k)fluoranthene	23.436	252	13581	0.389	ng/u1#	93
26) Benzo(a)pyrene	24.044	252	11196	0.410	ng/u1#	91
27) Indeno(1,2,3-cd)pyrene	26.801	276	13362	0.435	ng/u1#	97
28) Dibenzo(a,h)anthracene	26.824	278	11403	0.438	ng/u1	96
29) Benzo(g,h,i)perylene	27.619	276	12441	0.440	ng/u1	98

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

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