

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN12222\
 Data File : BN023363.D
 Acq On : 23 Dec 2022 06:25
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4353

Quant Time: Dec 23 06:56:17 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN122122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 22 02:28:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.994	152	7772	0.400	ng/ul	0.00
4) Naphthalene-d8	10.810	136	25619	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.645	164	15370	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.390	188	35397	0.400	ng/ul	0.00
17) Chrysene-d12	21.581	240	26507	0.400	ng/ul	0.00
23) Perylene-d12	24.030	264	20697	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.349	96	3300	0.391	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.399	152	16171	0.408	ng/ul	0.00
18) Fluoranthene-d10	19.420	212	39063	0.407	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.382	88	3167	0.367	ng/ul	93
5) Naphthalene	10.859	128	29631	0.394	ng/ul	100
7) 2-Methylnaphthalene	12.470	142	19333	0.402	ng/ul	100
8) 1-Methylnaphthalene	12.690	142	20023	0.404	ng/ul	100
10) Acenaphthylene	14.363	152	27520	0.404	ng/ul	100
11) Acenaphthene	14.705	153	22456	0.391	ng/ul	99
12) Fluorene	15.691	166	25843	0.401	ng/ul	99
14) Pentachlorophenol	17.040	266	4264	0.434	ng/ul	99
15) Phenanthrene	17.432	178	44398	0.385	ng/ul	100
16) Anthracene	17.525	178	38947	0.407	ng/ul	99
19) Fluoranthene	19.448	202	53525	0.416	ng/ul	100
20) Pyrene	19.811	202	54218	0.418	ng/ul	100
21) Benzo(a)anthracene	21.560	228	45765	0.443	ng/ul	100
22) Chrysene	21.616	228	44736	0.426	ng/ul	100
24) Benzo(b)fluoranthene	23.276	252	40389	0.377	ng/ul	99
25) Benzo(k)fluoranthene	23.326	252	38292	0.381	ng/ul	99
26) Benzo(a)pyrene	23.916	252	32670	0.385	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.590	276	38259	0.387	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.610	278	30075	0.392	ng/ul	99
29) Benzo(g,h,i)perylene	27.378	276	31296	0.388	ng/ul	100

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

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