

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122422\
 Data File : BN023393.D
 Acq On : 24 Dec 2022 04:41
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4356

Quant Time: Dec 24 05:12:32 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	7.990	152	7037	0.400	ng/u1	0.00
4) Naphthalene-d8	10.810	136	23258	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.641	164	13140	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.386	188	28339	0.400	ng/u1	0.00
17) Chrysene-d12	21.575	240	18930	0.400	ng/u1	0.00
23) Perylene-d12	24.025	264	14348	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.345	96	2811	0.368	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.394	152	14524	0.404	ng/u1	-0.01
18) Fluoranthene-d10	19.416	212	29960	0.438	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.378	88	2996	0.383	ng/u1	96
5) Naphthalene	10.859	128	27427	0.402	ng/u1	99
7) 2-Methylnaphthalene	12.471	142	17567	0.402	ng/u1	100
8) 1-Methylnaphthalene	12.685	142	18147	0.403	ng/u1	100
10) Acenaphthylene	14.358	152	24225	0.416	ng/u1	100
11) Acenaphthene	14.701	153	19592	0.399	ng/u1	99
12) Fluorene	15.686	166	21868	0.396	ng/u1	100
14) Pentachlorophenol	17.036	266	3248	0.413	ng/u1	99
15) Phenanthrene	17.428	178	35648	0.386	ng/u1	100
16) Anthracene	17.521	178	30825	0.403	ng/u1	100
19) Fluoranthene	19.448	202	39319	0.428	ng/u1	98
20) Pyrene	19.811	202	39497	0.427	ng/u1	98
21) Benzo(a)anthracene	21.560	228	31368	0.426	ng/u1	100
22) Chrysene	21.613	228	30474	0.407	ng/u1	100
24) Benzo(b)fluoranthene	23.273	252	26781	0.361	ng/u1	98
25) Benzo(k)fluoranthene	23.323	252	25324	0.363	ng/u1	99
26) Benzo(a)pyrene	23.911	252	21831	0.372	ng/u1	97
27) Indeno(1,2,3-cd)pyrene	26.580	276	25685	0.375	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.600	278	20248	0.381	ng/u1	98
29) Benzo(g,h,i)perylene	27.369	276	21315	0.381	ng/u1	99

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

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