

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081022\
 Data File : BN021124.D
 Acq On : 10 Aug 2022 19:02
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4280

Quant Time: Aug 10 22:09:00 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 09 23:10:06 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	2682	0.400	ng/ul	0.01
4) Naphthalene-d8	10.585	136	9206	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.438	164	5155	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.189	188	9659	0.400	ng/ul	0.00
17) Chrysene-d12	21.395	240	7084	0.400	ng/ul	0.00
23) Perylene-d12	23.725	264	5393	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.198	96	1276	0.372	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.186	152	5374	0.398	ng/ul	0.00
18) Fluoranthene-d10	19.226	212	9961	0.455	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.232	88	1440	0.407	ng/ul#	85
5) Naphthalene	10.635	128	9664	0.342	ng/ul	99
7) 2-Methylnaphthalene	12.263	142	5911	0.345	ng/ul	99
8) 1-Methylnaphthalene	12.477	142	6529	0.380	ng/ul	100
10) Acenaphthylene	14.161	152	8360	0.347	ng/ul	99
11) Acenaphthene	14.503	153	6750	0.339	ng/ul	99
12) Fluorene	15.493	166	7259	0.333	ng/ul	99
14) Pentachlorophenol	16.864	266	438	0.294	ng/ul	99
15) Phenanthrene	17.231	178	11342	0.328	ng/ul	100
16) Anthracene	17.328	178	9127	0.341	ng/ul	99
19) Fluoranthene	19.259	202	11928	0.380	ng/ul	99
20) Pyrene	19.621	202	12376	0.382	ng/ul	99
21) Benzo(a)anthracene	21.378	228	7867	0.340	ng/ul	100
22) Chrysene	21.430	228	9647	0.317	ng/ul	100
24) Benzo(b)fluoranthene	23.017	252	8141	0.326	ng/ul	95
25) Benzo(k)fluoranthene	23.064	252	8643	0.328	ng/ul	99
26) Benzo(a)pyrene	23.622	252	7576	0.333	ng/ul#	88
27) Indeno(1,2,3-cd)pyrene	26.129	276	7909	0.291	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.149	278	6208	0.289	ng/ul	97
29) Benzo(g,h,i)perylene	26.867	276	7002	0.289	ng/ul	97

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

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