

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081422\
 Data File : BN021224.D
 Acq On : 14 Aug 2022 19:16
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4289

Quant Time: Aug 15 00:42:55 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 : Sat Aug 13 04:02:44 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	2899	0.400	ng/ul	0.00
4) Naphthalene-d8	10.578	136	10229	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.432	164	5920	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.187	188	10832	0.400	ng/ul	0.00
17) Chrysene-d12	21.392	240	7341	0.400	ng/ul	0.00
23) Perylene-d12	23.719	264	5645	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.194	96	1445	0.390	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.184	152	5973	0.399	ng/ul	0.00
18) Fluoranthene-d10	19.225	212	11195	0.494	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.228	88	1485	0.389	ng/ul#	89
5) Naphthalene	10.628	128	11035	0.351	ng/ul	99
7) 2-Methylnaphthalene	12.256	142	6725	0.353	ng/ul	99
8) 1-Methylnaphthalene	12.476	142	7486	0.393	ng/ul	100
10) Acenaphthylene	14.155	152	9687	0.350	ng/ul	99
11) Acenaphthene	14.497	153	7839	0.342	ng/ul	98
12) Fluorene	15.492	166	8443	0.337	ng/ul	99
14) Pentachlorophenol	16.858	266	418	0.250	ng/ul	98
15) Phenanthrene	17.229	178	13036	0.336	ng/ul	99
16) Anthracene	17.326	178	10447	0.348	ng/ul	99
19) Fluoranthene	19.257	202	13630	0.420	ng/ul	100
20) Pyrene	19.620	202	13969	0.416	ng/ul	97
21) Benzo(a)anthracene	21.378	228	7877	0.329	ng/ul	99
22) Chrysene	21.427	228	10068	0.319	ng/ul	100
24) Benzo(b)fluoranthene	23.011	252	8447	0.323	ng/ul	96
25) Benzo(k)fluoranthene	23.058	252	8909	0.323	ng/ul	99
26) Benzo(a)pyrene	23.617	252	8071	0.339	ng/ul#	88
27) Indeno(1,2,3-cd)pyrene	26.132	276	8471	0.298	ng/ul	98
28) Dibenzo(a,h)anthracene	26.146	278	6516	0.290	ng/ul	97
29) Benzo(g,h,i)perylene	26.863	276	7762	0.306	ng/ul	98

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

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