

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062122\
 Data File : BN020334.D
 Acq On : 22 Jun 2022 08:15
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4376

Quant Time: Jun 22 08:46:44 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN061822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 20 01:13:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.800	152	727	0.400	ng/ul	0.00
4) Naphthalene-d8	10.579	136	2653	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.419	164	2115	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.163	188	5699	0.400	ng/ul	0.00
17) Chrysene-d12	21.348	240	8953	0.400	ng/ul	0.00
23) Perylene-d12	23.654	264	10797	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.260	96	332	0.400	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.173	152	1729	0.394	ng/ul	0.00
18) Fluoranthene-d10	19.189	212	7950	0.430	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.294	88	294	0.368	ng/ul#	87
5) Naphthalene	10.634	128	2932	0.346	ng/ul	99
7) 2-Methylnaphthalene	12.245	142	2013	0.367	ng/ul	99
8) 1-Methylnaphthalene	12.465	142	2173	0.402	ng/ul	99
10) Acenaphthylene	14.137	152	3872	0.383	ng/ul	99
11) Acenaphthene	14.479	153	3040	0.314	ng/ul	99
12) Fluorene	15.469	166	3959	0.395	ng/ul	98
14) Pentachlorophenol	16.829	266	662	0.605	ng/ul	99
15) Phenanthrene	17.205	178	7932	0.385	ng/ul	99
16) Anthracene	17.294	178	6967	0.399	ng/ul	99
19) Fluoranthene	19.221	202	10612	0.294	ng/ul	100
20) Pyrene	19.584	202	11565	0.295	ng/ul	96
21) Benzo(a)anthracene	21.333	228	11554	0.332	ng/ul	99
22) Chrysene	21.386	228	13482	0.312	ng/ul	98
24) Benzo(b)fluoranthene	22.955	252	16107	0.209	ng/ul	99
25) Benzo(k)fluoranthene	23.002	252	15818	0.220	ng/ul	99
26) Benzo(a)pyrene	23.558	252	16395	0.336	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.024	276	20683	0.260	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.038	278	16676	0.263	ng/ul	99
29) Benzo(g,h,i)perylene	26.745	276	18474	0.263	ng/ul	99

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

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