

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062623\
 Data File : BN026028.D
 Acq On : 26 Jun 2023 15:46
 Operator : MA/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4315

Quant Time: Jun 27 01:38:48 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 22 01:27:41 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.980	152	6003	0.400	ng/u1	0.00
4) Naphthalene-d8	10.787	136	20143	0.400	ng/u1	-0.01
9) Acenaphthene-d10	14.612	164	12194	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.359	188	24549	0.400	ng/u1	0.00
17) Chrysene-d12	21.546	240	14825	0.400	ng/u1	0.00
23) Perylene-d12	24.013	264	13563	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.331	96	2465	0.365	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.376	152	11958	0.386	ng/u1	0.00
18) Fluoranthene-d10	19.387	212	24967	0.442	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.364	88	2743	0.382	ng/u1	97
5) Naphthalene	10.836	128	22502	0.390	ng/u1	99
7) 2-Methylnaphthalene	12.447	142	14132	0.384	ng/u1	100
8) 1-Methylnaphthalene	12.667	142	15188	0.399	ng/u1	100
10) Acenaphthylene	14.334	152	23269	0.408	ng/u1	100
11) Acenaphthene	14.676	153	18330	0.392	ng/u1	99
12) Fluorene	15.662	166	20154	0.376	ng/u1	100
14) Pentachlorophenol	17.017	266	2350	0.267	ng/u1	99
15) Phenanthrene	17.402	178	31430	0.388	ng/u1	100
16) Anthracene	17.494	178	27974	0.388	ng/u1	99
19) Fluoranthene	19.414	202	34483	0.435	ng/u1	99
20) Pyrene	19.777	202	35206	0.427	ng/u1	98
21) Benzo(a)anthracene	21.529	228	23669	0.387	ng/u1	99
22) Chrysene	21.584	228	24635	0.387	ng/u1	99
24) Benzo(b)fluoranthene	23.256	252	21909	0.378	ng/u1	96
25) Benzo(k)fluoranthene	23.306	252	22389	0.385	ng/u1	97
26) Benzo(a)pyrene	23.905	252	19585	0.390	ng/u1	97
27) Indeno(1,2,3-cd)pyrene	26.598	276	24302	0.443	ng/u1#	96
28) Dibenzo(a,h)anthracene	26.621	278	18506	0.445	ng/u1	97
29) Benzo(g,h,i)perylene	27.389	276	20687	0.423	ng/u1	98

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

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