

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031824\
 Data File : BN030430.D
 Acq On : 18 Mar 2024 14:50
 Operator : MA/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4362

Quant Time: Mar 18 15:29:07 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN031524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 15 01:04:41 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.603	152	1906	0.400	ng/ul	0.00
4) Naphthalene-d8	10.376	136	4598	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.252	164	2779	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.011	188	5619	0.400	ng/ul	0.00
17) Chrysene-d12	21.220	240	5131	0.400	ng/ul	0.00
23) Perylene-d12	23.432	264	5863	0.400	ng/ul	# 0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.151	96	1023	0.440	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.981	152	2539	0.400	ng/ul	0.00
18) Fluoranthene-d10	19.049	212	6285	0.398	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.185	88	1105	0.437	ng/ul	94
5) Naphthalene	10.425	128	4848	0.391	ng/ul	98
7) 2-Methylnaphthalene	12.053	142	2896	0.398	ng/ul#	100
8) 1-Methylnaphthalene	12.273	142	3103	0.402	ng/ul#	28
10) Acenaphthylene	13.970	152	5129	0.387	ng/ul	99
11) Acenaphthene	14.317	153	3763	0.388	ng/ul	99
12) Fluorene	15.316	166	4256	0.398	ng/ul	100
14) Pentachlorophenol	16.665	266	666	0.412	ng/ul#	100
15) Phenanthrene	17.053	178	6524	0.394	ng/ul	99
16) Anthracene	17.150	178	5835	0.386	ng/ul	99
19) Fluoranthene	19.077	202	8060	0.397	ng/ul	99
20) Pyrene	19.444	202	8660	0.399	ng/ul	98
21) Benzo(a)anthracene	21.205	228	6696	0.389	ng/ul	99
22) Chrysene	21.255	228	7693	0.389	ng/ul	98
24) Benzo(b)fluoranthene	22.768	252	7733	0.393	ng/ul	96
25) Benzo(k)fluoranthene	22.812	252	9382	0.392	ng/ul	98
26) Benzo(a)pyrene	23.333	252	8072	0.389	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	25.645	276	10094	0.368	ng/ul#	87
28) Dibenzo(a,h)anthracene	25.672	278	7578	0.362	ng/ul	96
29) Benzo(g,h,i)perylene	26.320	276	9132	0.373	ng/ul	99

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

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