

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031524\
 Data File : BN030423.D
 Acq On : 15 Mar 2024 22:26
 Operator : MA/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4360

Quant Time: Mar 15 23:02:35 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	2082	0.400	ng/ul	0.00
4) Naphthalene-d8	10.380	136	4816	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.251	164	2851	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.010	188	5837	0.400	ng/ul	0.00
17) Chrysene-d12	21.214	240	5544	0.400	ng/ul	0.00
23) Perylene-d12	23.424	264	6454	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.155	96	1035	0.408	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.980	152	2648	0.398	ng/ul	0.00
18) Fluoranthene-d10	19.048	212	6569	0.385	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.189	88	1066	0.386	ng/ul	90
5) Naphthalene	10.429	128	4952	0.381	ng/ul	99
7) 2-Methylnaphthalene	12.057	142	3032	0.398	ng/ul#	100
8) 1-Methylnaphthalene	12.277	142	3248	0.402	ng/ul#	29
10) Acenaphthylene	13.974	152	5251	0.386	ng/ul	98
11) Acenaphthene	14.316	153	3870	0.389	ng/ul	100
12) Fluorene	15.315	166	4389	0.401	ng/ul	98
14) Pentachlorophenol	16.664	266	690	0.411	ng/ul#	100
15) Phenanthrene	17.052	178	6774	0.394	ng/ul	99
16) Anthracene	17.145	178	6077	0.387	ng/ul	98
19) Fluoranthene	19.076	202	8363	0.381	ng/ul	99
20) Pyrene	19.443	202	8993	0.384	ng/ul	97
21) Benzo(a)anthracene	21.200	228	7246	0.390	ng/ul	99
22) Chrysene	21.252	228	8386	0.392	ng/ul	98
24) Benzo(b)fluoranthene	22.760	252	8568	0.396	ng/ul	94
25) Benzo(k)fluoranthene	22.807	252	10154	0.386	ng/ul	95
26) Benzo(a)pyrene	23.327	252	8803	0.385	ng/ul	91
27) Indeno(1,2,3-cd)pyrene	25.636	276	11592	0.384	ng/ul#	88
28) Dibenzo(a,h)anthracene	25.656	278	8902	0.386	ng/ul	92
29) Benzo(g,h,i)perylene	26.310	276	10404	0.386	ng/ul	98

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

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