

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071323\
 Data File : BN026491.D
 Acq On : 13 Jul 2023 15:17
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4354

Quant Time: Jul 14 04:31:14 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN071223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 13 16:59:14 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.988	152	4249	0.400	ng/ul	0.00
4) Naphthalene-d8	10.792	136	14997	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.616	164	7792	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.368	188	12331	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.564	240	8604	0.400	ng/ul	0.00
23) Perylene-d12	24.002	264	9468	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.335	96	2114	0.378	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.381	152	6917	0.339	ng/ul	0.00
18) Fluoranthene-d10	19.396	212	11605	0.389	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.373	88	2000	0.378	ng/ul	94
5) Naphthalene	10.842	128	16263	0.395	ng/ul	99
7) 2-Methylnaphthalene	12.453	142	8035	0.344	ng/ul	100
8) 1-Methylnaphthalene	12.667	142	10065	0.402	ng/ul	99
10) Acenaphthylene	14.339	152	14144	0.371	ng/ul	98
11) Acenaphthene	14.676	153	11412	0.393	ng/ul	99
12) Fluorene	15.666	166	10565	0.346	ng/ul	99
14) Pentachlorophenol	17.030	266	1048	0.337	ng/ul	96
15) Phenanthrene	17.406	178	15243	0.388	ng/ul	99
16) Anthracene	17.503	178	12829	0.376	ng/ul	98
19) Fluoranthene	19.424	202	16382	0.406	ng/ul#	94
20) Pyrene	19.791	202	17251	0.394	ng/ul#	93
21) Benzo(a)anthracene	21.546	228	13117	0.415	ng/ul	99
22) Chrysene	21.599	228	14097	0.392	ng/ul	100
24) Benzo(b)fluoranthene	23.253	252	14069	0.381	ng/ul	94
25) Benzo(k)fluoranthene	23.303	252	14857	0.384	ng/ul#	93
26) Benzo(a)pyrene	23.894	252	13131	0.400	ng/ul	93
27) Indeno(1,2,3-cd)pyrene	26.551	276	16542	0.384	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.564	278	12557	0.380	ng/ul	94
29) Benzo(g,h,i)perylene	27.332	276	13697	0.361	ng/ul	99

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

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