

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071223\
 Data File : BN026479.D
 Acq On : 12 Jul 2023 17:58
 Operator : MA/JU
 Sample : SSTD1.617
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.6234

Quant Time: Jul 13 04:17:34 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 04:15:36 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.005	152	3422	0.400	ng/ul	0.00
4) Naphthalene-d8	10.803	136	11122	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.622	164	5375	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.369	188	8425	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.555	240	5775	0.400	ng/ul	0.00
23) Perylene-d12	24.002	264	6029	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.343	96	6211	0.462	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.387	152	24887	1.655	ng/ul	-0.01
18) Fluoranthene-d10	19.396	212	29824	1.523	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.381	88	6102	0.706	ng/ul	88
5) Naphthalene	10.853	128	49085	1.616	ng/ul	96
7) 2-Methylnaphthalene	12.458	142	29162	1.651	ng/ul	100
8) 1-Methylnaphthalene	12.678	142	27344	1.435	ng/ul	99
10) Acenaphthylene	14.344	152	41119	1.605	ng/ul	98
11) Acenaphthene	14.682	153	32620	1.694	ng/ul	99
12) Fluorene	15.672	166	32897	1.643	ng/ul	98
14) Pentachlorophenol	17.031	266	3133	1.525	ng/ul	98
15) Phenanthrene	17.411	178	43501	1.647	ng/ul	98
16) Anthracene	17.504	178	38622	1.601	ng/ul	96
19) Fluoranthene	19.424	202	41745	1.639	ng/ul	99
20) Pyrene	19.791	202	49095	1.773	ng/ul#	96
21) Benzo(a)anthracene	21.541	228	35638	1.725	ng/ul	97
22) Chrysene	21.593	228	39380	1.708	ng/ul	99
24) Benzo(b)fluoranthene	23.245	252	40131	1.748	ng/ul	85
25) Benzo(k)fluoranthene	23.298	252	42256	1.788	ng/ul#	86
26) Benzo(a)pyrene	23.888	252	35086	1.738	ng/ul#	79
27) Indeno(1,2,3-cd)pyrene	26.541	276	39817	1.562	ng/ul#	85
28) Dibenzo(a,h)anthracene	26.554	278	31071	1.565	ng/ul#	80
29) Benzo(g,h,i)perylene	27.326	276	34133	1.492	ng/ul#	88

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

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