

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101921\
 Data File : BN017003.D
 Acq On : 19 Oct 2021 13:13
 Operator : CG/JU
 Sample : SSTD1.626
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.6226

Quant Time: Oct 19 13:53:47 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN101921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 19 12:51:41 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.548	152	1894	0.40	ng/ul	0.00
4) Naphthalene-d8	10.322	136	7499	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.199	164	4720	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.953	188	10496	0.40	ng/ul	0.00
17) Chrysene-d12	21.162	240	12079	0.40	ng/ul	0.00
23) Perylene-d12	23.375	264	12019	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.076	96	3237	1.47	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.923	152	17001	1.61	ng/ul	0.00
18) Fluoranthene-d10	18.995	212	45055	1.39	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.109	88	3625	1.50	ng/ul#	54
5) Naphthalene	10.372	128	32724	1.64	ng/ul	98
7) 2-Methylnaphthalene	11.994	142	22683	1.74	ng/ul	100
8) 1-Methylnaphthalene	12.220	142	22814	1.75	ng/ul	100
10) Acenaphthylene	13.916	152	29910	1.68	ng/ul	98
11) Acenaphthene	14.263	153	25450	1.70	ng/ul	99
12) Fluorene	15.253	166	29991	1.75	ng/ul	99
14) Pentachlorophenol	16.607	266	3870	2.28	ng/ul	99
15) Phenanthrene	16.995	178	51913	1.72	ng/ul	98
16) Anthracene	17.084	178	45236	1.73	ng/ul	99
19) Fluoranthene	19.023	202	64185	1.60	ng/ul	97
20) Pyrene	19.390	202	63638	1.56	ng/ul	97
21) Benzo(a)anthracene	21.148	228	62079	1.68	ng/ul	99
22) Chrysene	21.200	228	68189	1.63	ng/ul	99
24) Benzo(b)fluoranthene	22.711	252	73935	1.63	ng/ul	97
25) Benzo(k)fluoranthene	22.755	252	78148	1.72	ng/ul	97
26) Benzo(a)pyrene	23.275	252	64421	1.68	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.603	276	95725	1.93	ng/ul	100
28) Dibenzo(a,h)anthracene	25.623	278	77026	1.89	ng/ul	98
29) Benzo(g,h,i)perylene	26.277	276	82134	1.91	ng/ul	98

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

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