

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102221\
 Data File : BN017063.D
 Acq On : 22 Oct 2021 01:47
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4276

Quant Time: Oct 22 03:10:57 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 : Thu Oct 21 10:28:56 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.548	152	1657	0.400	ng/ul	0.00
4) Naphthalene-d8	10.321	136	6368	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.197	164	3786	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.952	188	8572	0.400	ng/ul	0.00
17) Chrysene-d12	21.161	240	9690	0.400	ng/ul	0.00
23) Perylene-d12	23.370	264	12042	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.076	96	582	0.323	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.921	152	3335	0.367	ng/ul	0.00
18) Fluoranthene-d10	18.989	212	8618	0.360	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.109	88	569	0.263	ng/ul#	65
5) Naphthalene	10.365	128	6831	0.354	ng/ul	99
7) 2-Methylnaphthalene	11.998	142	4505	0.375	ng/ul	99
8) 1-Methylnaphthalene	12.218	142	4679	0.383	ng/ul	100
10) Acenaphthylene	13.915	152	6292	0.411	ng/ul	100
11) Acenaphthene	14.257	153	5179	0.394	ng/ul	97
12) Fluorene	15.256	166	5798	0.382	ng/ul	100
14) Pentachlorophenol	16.610	266	606	0.293	ng/ul	100
15) Phenanthrene	16.994	178	10733	0.394	ng/ul	100
16) Anthracene	17.082	178	8652	0.377	ng/ul	99
19) Fluoranthene	19.021	202	12430	0.374	ng/ul	99
20) Pyrene	19.384	202	12869	0.377	ng/ul	99
21) Benzo(a)anthracene	21.146	228	11379	0.360	ng/ul	99
22) Chrysene	21.196	228	14527	0.404	ng/ul	99
24) Benzo(b)fluoranthene	22.707	252	15810	0.349	ng/ul	100
25) Benzo(k)fluoranthene	22.753	252	19403	0.388	ng/ul	98
26) Benzo(a)pyrene	23.274	252	14769	0.365	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.598	276	23718	0.412	ng/ul	98
28) Dibenzo(a,h)anthracene	25.622	278	18307	0.400	ng/ul	98
29) Benzo(g,h,i)perylene	26.276	276	19490	0.391	ng/ul	100

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

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