

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070821\
 Data File : BN015412.D
 Acq On : 08 Jul 2021 17:09
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4111

Quant Time: Jul 08 17:50:52 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 08 11:17:31 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	3020	0.400	ng/ul	0.00
4) Naphthalene-d8	10.485	136	10293	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.340	164	5328	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.086	188	10747	0.400	ng/ul	0.00
17) Chrysene-d12	21.284	240	8603	0.400	ng/ul	0.00
23) Perylene-d12	23.523	264	7939	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	1307	0.384	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.074	152	5911	0.366	ng/ul	0.00
18) Fluoranthene-d10	19.118	212	10840	0.383	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.315	88	1374	0.374	ng/ul#	94
5) Naphthalene	10.529	128	12126	0.369	ng/ul	99
7) 2-Methylnaphthalene	12.146	142	7907	0.368	ng/ul	99
8) 1-Methylnaphthalene	12.371	142	7709	0.371	ng/ul	99
10) Acenaphthylene	14.057	152	9695	0.352	ng/ul	100
11) Acenaphthene	14.400	153	7789	0.369	ng/ul	99
12) Fluorene	15.390	166	8694	0.373	ng/ul	99
14) Pentachlorophenol	16.740	266	899	0.378	ng/ul	99
15) Phenanthrene	17.128	178	14331	0.367	ng/ul	99
16) Anthracene	17.217	178	12241	0.356	ng/ul	100
19) Fluoranthene	19.150	202	14887	0.374	ng/ul	100
20) Pyrene	19.513	202	15026	0.377	ng/ul	98
21) Benzo(a)anthracene	21.270	228	12341	0.387	ng/ul	99
22) Chrysene	21.322	228	13661	0.358	ng/ul	99
24) Benzo(b)fluoranthene	22.854	252	13834	0.373	ng/ul	94
25) Benzo(k)fluoranthene	22.898	252	14412	0.344	ng/ul	98
26) Benzo(a)pyrene	23.430	252	12551	0.403	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	25.774	276	15841	0.421	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.784	278	12572	0.426	ng/ul	99
29) Benzo(g,h,i)perylene	26.458	276	13336	0.383	ng/ul	98

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

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