

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111321\
 Data File : BN017458.D
 Acq On : 15 Nov 2021 09:24
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 66 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4294

Quant Time: Nov 15 12:57:09 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN111321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 15 02:31:43 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.864	152	7896	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.658	136	28303	0.400	ng/ul	#-0.02
9) Acenaphthene-d10	14.495	164	16031	0.400	ng/ul	-0.01
13) Phenanthrene-d10	17.236	188	31883	0.400	ng/ul	-0.02
17) Chrysene-d12	21.422	240	21184	0.400	ng/ul	-0.02
23) Perylene-d12	23.793	264	14228	0.400	ng/ul	#-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.328	96	3371	0.389	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.247	152	16385	0.375	ng/ul	-0.02
18) Fluoranthene-d10	19.264	212	31130	0.409	ng/ul	-0.02
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.362	88	4444	0.398	ng/ul#	39
5) Naphthalene	10.707	128	31355	0.401	ng/ul	96
7) 2-Methylnaphthalene	12.324	142	21445	0.383	ng/ul	100
8) 1-Methylnaphthalene	12.539	142	21180	0.382	ng/ul	99
10) Acenaphthylene	14.212	152	26888	0.349	ng/ul	99
11) Acenaphthene	14.559	153	22162	0.403	ng/ul	100
12) Fluorene	15.545	166	25170	0.386	ng/ul	98
14) Pentachlorophenol	16.890	266	2434	0.349	ng/ul	98
15) Phenanthrene	17.278	178	39951	0.417	ng/ul	99
16) Anthracene	17.371	178	35047	0.363	ng/ul	99
19) Fluoranthene	19.297	202	41795	0.425	ng/ul	98
20) Pyrene	19.659	202	41938	0.417	ng/ul	97
21) Benzo(a)anthracene	21.408	228	30687	0.390	ng/ul	99
22) Chrysene	21.460	228	32389	0.411	ng/ul	100
24) Benzo(b)fluoranthene	23.071	252	26325	0.453	ng/ul	93
25) Benzo(k)fluoranthene	23.120	252	26890	0.449	ng/ul#	92
26) Benzo(a)pyrene	23.688	252	20771	0.428	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	26.244	276	19427	0.474	ng/ul#	88
28) Dibenzo(a,h)anthracene	26.267	278	14428	0.473	ng/ul	93
29) Benzo(g,h,i)perylene	26.988	276	18788	0.501	ng/ul	96

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

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