

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121621\
 Data File : BN018099.D
 Acq On : 17 Dec 2021 14:37
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 75 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4301

Quant Time: Dec 17 17:59:23 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN121621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 15 17:47:27 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.662	152	4165	0.400	ng/ul	0.00
4) Naphthalene-d8	10.464	136	12328	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.313	164	7269	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.070	188	13427	0.400	ng/ul	0.00
17) Chrysene-d12	21.269	240	13938	0.400	ng/ul #	0.00
23) Perylene-d12	23.546	264	11884	0.400	ng/ul #	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.139	96	1790	0.414	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.070	152	5483	0.382	ng/ul	0.00
18) Fluoranthene-d10	19.105	212	16192	0.408	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.172	88	2051	0.404	ng/ul#	46
5) Naphthalene	10.513	128	12897	0.382	ng/ul	99
7) 2-Methylnaphthalene	12.141	142	7129	0.375	ng/ul	100
8) 1-Methylnaphthalene	12.350	142	8256	0.374	ng/ul	100
10) Acenaphthylene	14.035	152	11097	0.377	ng/ul	99
11) Acenaphthene	14.378	153	9994	0.382	ng/ul	97
12) Fluorene	15.377	166	10737	0.372	ng/ul	99
14) Pentachlorophenol	16.749	266	714	0.304	ng/ul	98
15) Phenanthrene	17.112	178	17891	0.384	ng/ul	99
16) Anthracene	17.209	178	13642	0.372	ng/ul	99
19) Fluoranthene	19.133	202	22561	0.402	ng/ul	98
20) Pyrene	19.495	202	25109	0.399	ng/ul#	94
21) Benzo(a)anthracene	21.254	228	15115	0.365	ng/ul	98
22) Chrysene	21.304	228	22957	0.386	ng/ul	99
24) Benzo(b)fluoranthene	22.856	252	19176	0.404	ng/ul	87
25) Benzo(k)fluoranthene	22.900	252	24027	0.379	ng/ul#	83
26) Benzo(a)pyrene	23.449	252	17830	0.389	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	25.883	276	19693	0.381	ng/ul#	91
28) Dibenzo(a,h)anthracene	25.907	278	14304	0.379	ng/ul#	86
29) Benzo(g,h,i)perylene	26.588	276	17186	0.388	ng/ul#	89

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

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