

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082621\
 Data File : BN016098.D
 Acq On : 27 Aug 2021 09:47
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4261

Quant Time: Aug 27 10:25:02 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN082621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 26 11:35:43 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.867	152	1346	0.400	ng/ul	0.00
4) Naphthalene-d8	10.666	136	6061	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.501	164	3767	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.246	188	7886	0.400	ng/ul	0.00
17) Chrysene-d12	21.433	240	7915	0.400	ng/ul	0.00
23) Perylene-d12	23.821	264	7370	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	717	0.393	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.250	152	3794	0.416	ng/ul	0.00
18) Fluoranthene-d10	19.276	212	8700	0.408	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.352	88	713	0.387	ng/ul	98
5) Naphthalene	10.710	128	6658	0.425	ng/ul	99
7) 2-Methylnaphthalene	12.327	142	4603	0.435	ng/ul	98
8) 1-Methylnaphthalene	12.541	142	4285	0.413	ng/ul	99
10) Acenaphthylene	14.219	152	6154	0.463	ng/ul	98
11) Acenaphthene	14.561	153	4877	0.425	ng/ul	99
12) Fluorene	15.547	166	5547	0.408	ng/ul	97
14) Pentachlorophenol	16.896	266	735	0.445	ng/ul	99
15) Phenanthrene	17.284	178	9017	0.414	ng/ul	99
16) Anthracene	17.377	178	8284	0.441	ng/ul	99
19) Fluoranthene	19.303	202	10534	0.426	ng/ul	99
20) Pyrene	19.666	202	10720	0.431	ng/ul	99
21) Benzo(a)anthracene	21.415	228	10750	0.436	ng/ul	99
22) Chrysene	21.468	228	11016	0.415	ng/ul	100
24) Benzo(b)fluoranthene	23.093	252	11024	0.401	ng/ul	98
25) Benzo(k)fluoranthene	23.140	252	10847	0.391	ng/ul	98
26) Benzo(a)pyrene	23.716	252	9688	0.419	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.286	276	12142	0.397	ng/ul#	58
28) Dibenzo(a,h)anthracene	26.306	278	10103	0.392	ng/ul	99
29) Benzo(g,h,i)perylene	27.044	276	10310	0.391	ng/ul	98

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

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