

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071721\
 Data File : BN015609.D
 Acq On : 16 Jul 2021 23:41
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.4252

Quant Time: Jul 17 00:44:55 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 : Fri Jul 16 15:41:51 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.791	152	2677	0.400	ng/ul	# 0.00
4) Naphthalene-d8	10.578	136	9569	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.427	164	5291	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.179	188	10965	0.400	ng/ul	0.00
17) Chrysene-d12	21.371	240	10047	0.400	ng/ul	# 0.00
23) Perylene-d12	23.713	264	9477	0.400	ng/ul	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.289	96	1151	0.381	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.173	152	5593	0.373	ng/ul	0.00
18) Fluoranthene-d10	19.210	212	11636	0.352	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.327	88	1231	0.378	ng/ul#	78
5) Naphthalene	10.628	128	11142	0.365	ng/ul#	90
7) 2-Methylnaphthalene	12.245	142	7407	0.371	ng/ul	99
8) 1-Methylnaphthalene	12.465	142	7307	0.378	ng/ul	99
10) Acenaphthylene	14.145	152	9788	0.358	ng/ul#	90
11) Acenaphthene	14.492	153	7700	0.367	ng/ul	99
12) Fluorene	15.478	166	8584	0.371	ng/ul	98
14) Pentachlorophenol	16.833	266	855	0.352	ng/ul	92
15) Phenanthrene	17.221	178	14321	0.360	ng/ul	98
16) Anthracene	17.310	178	12631	0.360	ng/ul	95
19) Fluoranthene	19.238	202	15708	0.338	ng/ul#	90
20) Pyrene	19.605	202	16104	0.346	ng/ul#	91
21) Benzo(a)anthracene	21.357	228	14090	0.379	ng/ul	98
22) Chrysene	21.407	228	14773	0.332	ng/ul	99
24) Benzo(b)fluoranthene	23.002	252	14956	0.338	ng/ul	93
25) Benzo(k)fluoranthene	23.049	252	15897	0.318	ng/ul#	93
26) Benzo(a)pyrene	23.610	252	13355	0.359	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	26.125	276	17464	0.389	ng/ul#	48
28) Dibenzo(a,h)anthracene	26.145	278	13908	0.395	ng/ul#	88
29) Benzo(g,h,i)perylene	26.866	276	14150	0.340	ng/ul#	86

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

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