

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031722\
 Data File : BN018995.D
 Acq On : 17 Mar 2022 15:30
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDCCC0.4

Quant Time: Mar 17 16:06:37 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN031722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 17 13:18:23 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Christian Giraldo 03/18/2022
 Supervised By :Jagrut Upadhyay 03/18/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.854	152	4514	0.400	ng	0.00
7) Naphthalene-d8	10.658	136	12282	0.400	ng	# 0.00
13) Acenaphthene-d10	14.495	164	8026	0.400	ng	0.00
19) Phenanthrene-d10	17.236	188	17077	0.400	ng	# 0.00
29) Chrysene-d12	21.422	240	12366	0.400	ng	# 0.00
35) Perylene-d12	23.793	264	9147	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.406	112	3985	0.394	ng	0.00
5) Phenol-d6	7.009	99	3952	0.373	ng	0.00
8) Nitrobenzene-d5	9.014	82	3153	0.356	ng	0.00
11) 2-Methylnaphthalene-d10	12.250	152	7864	0.386	ng	0.00
14) 2,4,6-Tribromophenol	15.985	330	1327	0.343	ng	0.00
15) 2-Fluorobiphenyl	13.127	172	12398	0.385	ng	0.00
27) Fluoranthene-d10	19.265	212	19770	0.390	ng	0.00
31) Terphenyl-d14	19.864	244	11742	0.387	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.247	88	2298	0.412	ng	97
3) n-Nitrosodimethylamine	3.572	42	2395	0.368	ng	# 95
6) bis(2-Chloroethyl)ether	7.269	93	4896	0.388	ng	98
9) Naphthalene	10.712	128	13628	0.390	ng	99
10) Hexachlorobutadiene	11.011	225	3437	0.402	ng	# 100
12) 2-Methylnaphthalene	12.325	142	9281	0.391	ng	99
16) Acenaphthylene	14.207	152	13407	0.360	ng	100
17) Acenaphthene	14.559	154	10054	0.385	ng	99
18) Fluorene	15.543	166	12782	0.388	ng	100
20) 4,6-Dinitro-2-methylph...	15.618	198	786	0.354	ng	# 78
21) 4-Bromophenyl-phenylether	16.426	248	4134	0.384	ng	# 87
22) Hexachlorobenzene	16.549	284	5477	0.395	ng	99
23) Atrazine	16.692	200	2809	0.349	ng	96
24) Pentachlorophenol	16.897	266	1364	0.320	ng	99
25) Phenanthrene	17.277	178	21069	0.395	ng	100
26) Anthracene	17.369	178	16262	0.360	ng	100
28) Fluoranthene	19.295	202	24470	0.391	ng	99
30) Pyrene	19.657	202	24873	0.404	ng	99
32) Benzo(a)anthracene	21.404	228	13560	0.356	ng	99
33) Chrysene	21.458	228	21188	0.412	ng	99
34) Bis(2-ethylhexyl)phtha...	21.333	149	9159	0.363	ng	99
36) Indeno(1,2,3-cd)pyrene	26.252	276	12912m	0.391	ng	
37) Benzo(b)fluoranthene	23.071	252	12791m	0.383	ng	
38) Benzo(k)fluoranthene	23.121	252	17918	0.392	ng	97
39) Benzo(a)pyrene	23.694	252	11897	0.392	ng	# 80
40) Dibenzo(a,h)anthracene	26.273	278	10063m	0.391	ng	
41) Benzo(g,h,i)perylene	26.998	276	12024	0.371	ng	97

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

