

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111921\
 Data File : BN017516.D
 Acq On : 19 Nov 2021 10:29
 Operator : CG/JU
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDICC0.1

Quant Time: Nov 19 12:30:32 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN111921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 19 12:29:22 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.865	152	28211	0.400	ng	0.00
7) Naphthalene-d8	10.648	136	127464	0.400	ng	#-0.01
13) Acenaphthene-d10	14.485	164	67933	0.400	ng	0.00
19) Phenanthrene-d10	17.225	188	130892	0.400	ng	# 0.00
29) Chrysene-d12	21.409	240	112442	0.400	ng	# 0.00
35) Perylene-d12	23.776	264	97956	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.440	112	6856	0.102	ng	0.00
5) Phenol-d6	7.026	99	7604	0.105	ng	0.00
8) Nitrobenzene-d5	9.001	82	7610	0.096	ng	-0.01
11) 2-Methylnaphthalene-d10	12.239	152	20579	0.098	ng	0.00
14) 2,4,6-Tribromophenol	15.970	330	1863	0.108	ng	0.00
15) 2-Fluorobiphenyl	13.115	172	24742	0.097	ng	0.00
27) Fluoranthene-d10	19.258	212	36321	0.093	ng	0.00
31) Terphenyl-d14	19.858	244	24229	0.088	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.383	88	1348	0.109	ng	# 1
3) n-Nitrosodimethylamine	3.715	42	2537	0.107	ng	# 90
6) bis(2-Chloroethyl)ether	7.284	93	8160	0.105	ng	95
9) Naphthalene	10.699	128	32816	0.102	ng	99
10) Hexachlorobutadiene	11.003	225	6150	0.092	ng	# 99
12) 2-Methylnaphthalene	12.315	142	20492	0.101	ng	98
16) Acenaphthylene	14.206	152	22602	0.089	ng	100
17) Acenaphthene	14.549	154	17792	0.097	ng	99
18) Fluorene	15.526	166	21028	0.100	ng	99
20) 4,6-Dinitro-2-methylph...	15.615	198	263	0.051	ng	# 8
21) 4-Bromophenyl-phenylether	16.422	248	6511	0.088	ng	# 82
22) Hexachlorobenzene	16.544	284	7464	0.089	ng	95
23) Atrazine	16.690	200	4070	0.096	ng	# 96
24) Pentachlorophenol	16.885	266	1728	0.103	ng	98
25) Phenanthrene	17.262	178	35144	0.098	ng	99
26) Anthracene	17.359	178	26832	0.090	ng	100
28) Fluoranthene	19.287	202	36716	0.095	ng	100
30) Pyrene	19.650	202	36998	0.092	ng	100
32) Benzo(a)anthracene	21.398	228	30881	0.089	ng	99
33) Chrysene	21.451	228	36576	0.102	ng	99
34) Bis(2-ethylhexyl)phtha...	21.335	149	12414	0.068	ng	100
36) Indeno(1,2,3-cd)pyrene	26.220	276	25087	0.111	ng	95
37) Benzo(b)fluoranthene	23.058	252	31163	0.085	ng	# 96
38) Benzo(k)fluoranthene	23.106	252	31899	0.094	ng	# 95
39) Benzo(a)pyrene	23.674	252	24882	0.087	ng	# 94
40) Dibenzo(a,h)anthracene	26.241	278	20472	0.119	ng	# 90
41) Benzo(g,h,i)perylene	26.967	276	21820	0.103	ng	96

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

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