

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102620\
 Data File : BN012393.D
 Acq On : 26 Oct 2020 10:51
 Operator : CG/JU
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDICC0.1

Quant Time: Oct 26 12:47:51 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN102620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 26 12:39:15 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.586	152	528	0.40	ng	# 0.00
7) Naphthalene-d8	10.364	136	2062	0.40	ng	# 0.01
13) Acenaphthene-d10	14.230	164	1236	0.40	ng	0.01
19) Phenanthrene-d10	16.980	188	2572	0.40	ng	# 0.01
29) Chrysene-d12	21.184	240	2692	0.40	ng	# 0.00
36) Perylene-d12	23.357	264	3204	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.194	112	205	0.14	ng	0.00
5) Phenol-d6	6.773	99	216	0.12	ng	0.00
8) Nitrobenzene-d5	8.755	82	192	0.09	ng	0.03
11) 2-Methylnaphthalene-d10	11.976	152	329	0.09	ng	0.02
14) 2,4,6-Tribromophenol	15.740	330	88	0.13	ng	0.01
15) 2-Fluorobiphenyl	12.860	172	437	0.08	ng	0.01
27) Fluoranthene-d10	19.023	212	847	0.10	ng	0.01
31) Terphenyl-d14	19.633	244	671	0.09	ng	0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.181	88	119	0.17	ng	# 84
6) bis(2-Chloroethyl)ether	7.031	93	119	0.08	ng	# 76
9) Naphthalene	10.402	128	634	0.11	ng	# 57
10) Hexachlorobutadiene	10.681	225	138	0.09	ng	# 99
12) 2-Methylnaphthalene	12.051	142	369	0.09	ng	# 80
16) Acenaphthylene	13.951	152	617	0.10	ng	97
17) Acenaphthene	14.294	154	405	0.09	ng	94
18) Fluorene	15.296	166	499	0.09	ng	94
20) 4,6-Dinitro-2-methylph...	15.435	198	48	0.08	ng	# 1
21) 4-Bromophenyl-phenylether	16.189	248	157	0.10	ng	# 86
22) Hexachlorobenzene	16.287	284	202	0.11	ng	# 84
23) Atrazine	16.457	200	153	0.10	ng	# 56
25) Phenanthrene	17.017	178	770	0.09	ng	96
26) Anthracene	17.127	178	684	0.09	ng	97
28) Fluoranthene	19.052	202	920	0.09	ng	# 98
30) Pyrene	19.415	202	995	0.10	ng	98
32) Benzo(a)anthracene	21.174	228	766	0.08	ng	# 82
33) Chrysene	21.216	228	1041	0.10	ng	# 87
34) Bis(2-ethylhexyl)phtha...	21.099	149	624	0.13	ng	# 90
35) Indeno(1,2,3-cd)pyrene	25.531	276	1435	0.11	ng	# 95
37) Benzo(b)fluoranthene	22.714	252	1034	0.08	ng	# 40
38) Benzo(k)fluoranthene	22.755	252	1241	0.09	ng	# 43
39) Benzo(a)pyrene	23.265	252	1027	0.09	ng	# 29
40) Dibenzo(a,h)anthracene	25.541	278	1147	0.09	ng	# 33
41) Benzo(g,h,i)perylene	26.178	276	1280	0.10	ng	# 75

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

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