

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092420\
 Data File : BN011911.D
 Acq On : 24 Sep 2020 13:31
 Operator : CG/JU
 Sample : L3988-22MS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 FC-MW03SR1-20200910MS

Manual Integrations
 APPROVED

mohammad
 9/25/2020 8:06:47 AM

Quant Time: Sep 24 14:05:00 2020
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN092320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 23 18:41:41 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.683	152	3690	0.40	ng	0.00
7) Naphthalene-d8	10.453	136	15746	0.40	ng	#-0.01
13) Acenaphthene-d10	14.319	164	8963m	0.40	ng	0.00
19) Phenanthrene-d10	17.054	188	19780	0.40	ng	#-0.01
29) Chrysene-d12	21.259	240	18019	0.40	ng	#-0.01
36) Perylene-d12	23.474	264	18053	0.40	ng	#-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.291	112	1154	0.09	ng	0.00
5) Phenol-d6	6.853	99	803	0.05	ng	0.00
8) Nitrobenzene-d5	8.818	82	3131	0.19	ng	-0.01
11) 2-Methylnaphthalene-d10	12.051	152	10397	0.42	ng	0.00
14) 2,4,6-Tribromophenol	15.803	330	1252	0.29	ng	-0.01
15) 2-Fluorobiphenyl	12.936	172	6422	0.19	ng	0.00
27) Fluoranthene-d10	19.097	212	25971	0.45	ng	0.00
31) Terphenyl-d14	19.703	244	13399	0.33	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.254	88	2159	0.34	ng	# 76
3) n-Nitrosodimethylamine	3.560	42	2189	0.24	ng	# 98
6) bis(2-Chloroethyl)ether	7.111	93	2904	0.21	ng	87
9) Naphthalene	10.504	128	349931	7.76	ng	98
10) Hexachlorobutadiene	10.795	225	1379	0.19	ng	# 99
12) 2-Methylnaphthalene	12.122	142	19126	0.67	ng	95
16) Acenaphthylene	14.027	152	10415	0.24	ng	# 82
17) Acenaphthene	14.370	154	8263	0.28	ng	# 89
18) Fluorene	15.359	166	13617	0.38	ng	# 96
20) 4,6-Dinitro-2-methylph...	15.448	198	959	0.25	ng	# 1
21) 4-Bromophenyl-phenylether	16.262	248	2114	0.20	ng	# 80
22) Hexachlorobenzene	16.372	284	2268	0.19	ng	# 86
23) Atrazine	16.530	200	1725	0.17	ng	# 84
24) Pentachlorophenol	16.713	266	2661	0.45	ng	97
25) Phenanthrene	17.102	178	14413	0.24	ng	97
26) Anthracene	17.187	178	11647	0.22	ng	98
28) Fluoranthene	19.127	202	13988	0.21	ng	# 94
30) Pyrene	19.489	202	14818	0.22	ng	98
32) Benzo(a)anthracene	21.248	228	12266	0.21	ng	98
33) Chrysene	21.291	228	12837	0.21	ng	98
34) Bis(2-ethylhexyl)phtha...	21.184	149	12563	0.31	ng	# 92
35) Indeno(1,2,3-cd)pyrene	25.695	276	14630	0.21	ng	# 89
37) Benzo(b)fluoranthene	22.810	252	13256	0.21	ng	# 69
38) Benzo(k)fluoranthene	22.854	252	13272	0.20	ng	# 77
39) Benzo(a)pyrene	23.378	252	11460	0.20	ng	# 44
40) Dibenzo(a,h)anthracene	25.705	278	11781	0.20	ng	# 80
41) Benzo(g,h,i)perylene	26.369	276	12628	0.20	ng	# 88

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

