

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080422\
 Data File : BN020984.D
 Acq On : 04 Aug 2022 11:16
 Operator : CG/JU
 Sample : N3900-08
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0C58

Quant Time: Aug 04 11:49:58 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 03 16:31:11 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	3177	0.400	ng/ul	0.00
4) Naphthalene-d8	10.602	136	10191	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.457	164	5766	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.210	188	10018	0.400	ng/ul	0.00
17) Chrysene-d12	21.418	240	9747	0.400	ng/ul	0.00
23) Perylene-d12	23.763	264	7733	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.207	96	2401	0.591	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.202	152	4872	0.326	ng/ul	0.00
18) Fluoranthene-d10	19.249	212	9545	0.317	ng/ul	0.00
Target Compounds					Qvalue	
5) Naphthalene	10.657	128	7007	0.224	ng/ul	99
7) 2-Methylnaphthalene	12.279	142	4615	0.243	ng/ul	98
10) Acenaphthylene	14.175	152	17303	0.642	ng/ul	98
11) Acenaphthene	14.517	153	6593	0.296	ng/ul	98
12) Fluorene	15.512	166	6722	0.276	ng/ul	99
15) Phenanthrene	17.248	178	21101	0.589	ng/ul	99
16) Anthracene	17.353	178	5753	0.207	ng/ul	98
19) Fluoranthene	19.277	202	13669	0.317	ng/ul	97
20) Pyrene	19.640	202	14496	0.326	ng/ul	97
21) Benzo(a)anthracene	21.404	228	21536	0.677	ng/ul	99
22) Chrysene	21.456	228	14828	0.354	ng/ul	99
24) Benzo(b)fluoranthene	23.055	252	12223	0.341	ng/ul	98
26) Benzo(a)pyrene	23.660	252	13453	0.412	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.186	276	12868	0.330	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.206	278	10612	0.344	ng/ul	97
29) Benzo(g,h,i)perylene	26.924	276	13799	0.397	ng/ul	98

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

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