

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080422\
 Data File : BN021019.D
 Acq On : 05 Aug 2022 10:19
 Operator : CG/JU
 Sample : PB146742BS
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS742

Quant Time: Aug 06 00:27:32 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	7.809	152	2226	0.400	ng/ul	0.00
4) Naphthalene-d8	10.602	136	7471	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.457	164	4237	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.205	188	8224	0.400	ng/ul	0.00
17) Chrysene-d12	21.415	240	6818	0.400	ng/ul	0.00
23) Perylene-d12	23.757	264	5751	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	2059	0.723	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.202	152	4326	0.395	ng/ul	0.00
18) Fluoranthene-d10	19.245	212	8825	0.419	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.240	88	5625	1.917	ng/ul#	83
5) Naphthalene	10.651	128	7829	0.341	ng/ul	99
7) 2-Methylnaphthalene	12.279	142	4777	0.343	ng/ul	99
8) 1-Methylnaphthalene	12.493	142	5251	0.377	ng/ul	100
10) Acenaphthylene	14.174	152	6863	0.346	ng/ul	100
11) Acenaphthene	14.517	153	5509	0.336	ng/ul	99
12) Fluorene	15.511	166	6001	0.335	ng/ul	99
14) Pentachlorophenol	16.872	266	891	0.703	ng/ul	98
15) Phenanthrene	17.248	178	9625	0.327	ng/ul	100
16) Anthracene	17.345	178	7871	0.345	ng/ul	99
19) Fluoranthene	19.277	202	10706	0.355	ng/ul	99
20) Pyrene	19.640	202	11057	0.355	ng/ul	97
21) Benzo(a)anthracene	21.401	228	8137	0.366	ng/ul	99
22) Chrysene	21.453	228	10207	0.348	ng/ul	99
24) Benzo(b)fluoranthene	23.043	252	9277	0.348	ng/ul	100
25) Benzo(k)fluoranthene	23.093	252	9590	0.341	ng/ul	100
26) Benzo(a)pyrene	23.651	252	8486	0.349	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.172	276	9090	0.313	ng/ul	100
28) Dibenzo(a,h)anthracene	26.196	278	7288	0.318	ng/ul	98
29) Benzo(g,h,i)perylene	26.913	276	8172	0.316	ng/ul	98

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

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