

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN030224\
 Data File : BN030222.D
 Acq On : 03 Mar 2024 16:59
 Operator : MA/JU
 Sample : PB159246BS
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS246

Quant Time: Mar 03 22:54:01 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN030224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 02 16:09:30 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	1778	0.400	ng/ul	0.00
4) Naphthalene-d8	10.628	136	4419	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.479	164	2082	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.239	188	3702	0.400	ng/ul	0.00
17) Chrysene-d12	21.447	240	3143	0.400	ng/ul	# 0.00
23) Perylene-d12	23.803	264	3707	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.307	96	1400	0.569	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.234	152	1700	0.298	ng/ul	0.00
18) Fluoranthene-d10	19.277	212	2981	0.291	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.336	88	3800	1.423	ng/ul#	87
5) Naphthalene	10.678	128	3626	0.278	ng/ul	98
7) 2-Methylnaphthalene	12.306	142	1981	0.280	ng/ul	100
8) 1-Methylnaphthalene	12.520	142	2125	0.284	ng/ul	99
10) Acenaphthylene	14.201	152	3233	0.278	ng/ul	98
11) Acenaphthene	14.544	153	2341	0.274	ng/ul	98
12) Fluorene	15.543	166	2397	0.275	ng/ul	98
14) Pentachlorophenol	16.893	266	810	0.767	ng/ul	98
15) Phenanthrene	17.281	178	3392	0.273	ng/ul	99
16) Anthracene	17.378	178	3091	0.283	ng/ul	99
19) Fluoranthene	19.305	202	3862	0.277	ng/ul	99
20) Pyrene	19.667	202	4141	0.273	ng/ul	97
21) Benzo(a)anthracene	21.432	228	3271	0.294	ng/ul	99
22) Chrysene	21.482	228	3791	0.284	ng/ul	98
24) Benzo(b)fluoranthene	23.087	252	4023	0.281	ng/ul	98
25) Benzo(k)fluoranthene	23.136	252	4511	0.268	ng/ul	97
26) Benzo(a)pyrene	23.701	252	4127	0.284	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.236	276	5414	0.291	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.262	278	4300	0.299	ng/ul	97
29) Benzo(g,h,i)perylene	26.970	276	4896	0.284	ng/ul#	94

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

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