

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN030224\
 Data File : BN030247.D
 Acq On : 04 Mar 2024 09:12
 Operator : MA/JU
 Sample : PB159339BS
 Misc :
 ALS Vial : 70 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS339

Quant Time: Mar 04 09:41:39 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 : Mon Mar 04 05:54:20 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	1143	0.400	ng/ul	0.00
4) Naphthalene-d8	10.628	136	3128	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.479	164	1497	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.239	188	2652	0.400	ng/ul	0.00
17) Chrysene-d12	21.447	240	2240	0.400	ng/ul	0.00
23) Perylene-d12	23.806	264	2586	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.307	96	1250	0.790	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.234	152	1778	0.440	ng/ul	0.00
18) Fluoranthene-d10	19.277	212	3150	0.432	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.336	88	3594	2.093	ng/ul#	87
5) Naphthalene	10.683	128	3846	0.417	ng/ul	100
7) 2-Methylnaphthalene	12.306	142	2133	0.426	ng/ul	99
8) 1-Methylnaphthalene	12.520	142	2270	0.428	ng/ul	99
10) Acenaphthylene	14.201	152	3527	0.421	ng/ul	97
11) Acenaphthene	14.544	153	2567	0.418	ng/ul	96
12) Fluorene	15.543	166	2599	0.415	ng/ul	97
14) Pentachlorophenol	16.888	266	843	1.114	ng/ul	98
15) Phenanthrene	17.281	178	3720	0.418	ng/ul	97
16) Anthracene	17.378	178	3303	0.422	ng/ul	98
19) Fluoranthene	19.305	202	4137	0.416	ng/ul	100
20) Pyrene	19.667	202	4484	0.415	ng/ul	96
21) Benzo(a)anthracene	21.435	228	3503	0.442	ng/ul	98
22) Chrysene	21.485	228	4033	0.424	ng/ul	99
24) Benzo(b)fluoranthene	23.087	252	4238	0.425	ng/ul	97
25) Benzo(k)fluoranthene	23.136	252	4902	0.417	ng/ul#	96
26) Benzo(a)pyrene	23.701	252	4322	0.426	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	26.239	276	5679	0.438	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.272	278	4432	0.442	ng/ul	97
29) Benzo(g,h,i)perylene	26.973	276	5084	0.423	ng/ul#	93

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

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