

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040124\
 Data File : BN030579.D
 Acq On : 31 Mar 2024 14:50
 Operator : MA/JU
 Sample : PB159893BS
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS893

Quant Time: Mar 31 15:50:48 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN032724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 27 15:11:42 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.704	152	7001	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.481	136	21519	0.400	ng/ul	-0.03
9) Acenaphthene-d10	14.341	164	12191	0.400	ng/ul	-0.02
13) Phenanthrene-d10	17.088	188	26863	0.400	ng/ul	-0.02
17) Chrysene-d12	21.281	240	23275	0.400	ng/ul	-0.02
23) Perylene-d12	23.517	264	21330	0.400	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.202	96	5068	0.661	ng/ul	-0.02
6) 2-Methylnaphthalene-d10	12.076	152	11288	0.348	ng/ul	-0.03
18) Fluoranthene-d10	19.120	212	25760	0.358	ng/ul	-0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.236	88	14300	1.840	ng/ul#	86
5) Naphthalene	10.530	128	21202	0.361	ng/ul	100
7) 2-Methylnaphthalene	12.152	142	13770	0.355	ng/ul	100
8) 1-Methylnaphthalene	12.372	142	14213	0.347	ng/ul	100
10) Acenaphthylene	14.059	152	20723	0.377	ng/ul	99
11) Acenaphthene	14.401	153	15720	0.367	ng/ul	99
12) Fluorene	15.391	166	17905	0.355	ng/ul	99
14) Pentachlorophenol	16.733	266	4049	0.824	ng/ul	100
15) Phenanthrene	17.130	178	30362	0.377	ng/ul	100
16) Anthracene	17.218	178	25978	0.383	ng/ul	99
19) Fluoranthene	19.152	202	35133	0.366	ng/ul	99
20) Pyrene	19.515	202	35908	0.380	ng/ul	100
21) Benzo(a)anthracene	21.264	228	32215	0.401	ng/ul	100
22) Chrysene	21.316	228	34302	0.387	ng/ul	99
24) Benzo(b)fluoranthene	22.842	252	32443	0.373	ng/ul	99
25) Benzo(k)fluoranthene	22.886	252	34574	0.378	ng/ul	100
26) Benzo(a)pyrene	23.421	252	28611	0.406	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.780	276	30869	0.387	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.800	278	24782	0.391	ng/ul	98
29) Benzo(g,h,i)perylene	26.468	276	24042	0.355	ng/ul	99

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

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