

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022824\
 Data File : BN030119.D
 Acq On : 28 Feb 2024 08:12
 Operator : MA/JU
 Sample : PB159258BS
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS258

Quant Time: Feb 28 08:38:34 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN020624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 23:49:29 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.851	152	3858	0.400	ng/ul	0.00
4) Naphthalene-d8	10.649	136	7858	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.496	164	4354	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.258	188	8733	0.400	ng/ul	0.00
17) Chrysene-d12	21.459	240	7031	0.400	ng/ul	0.00
23) Perylene-d12	23.820	264	5897	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	3933	0.818	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.244	152	4541	0.454	ng/ul	0.00
18) Fluoranthene-d10	19.289	212	9658	0.418	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.349	88	11155	2.130	ng/ul#	82
5) Naphthalene	10.699	128	8495	0.384	ng/ul	99
7) 2-Methylnaphthalene	12.315	142	5523	0.436	ng/ul	100
8) 1-Methylnaphthalene	12.535	142	5867	0.454	ng/ul	98
10) Acenaphthylene	14.218	152	8039	0.378	ng/ul	97
11) Acenaphthene	14.561	153	6445	0.399	ng/ul	99
12) Fluorene	15.560	166	7309	0.433	ng/ul	97
14) Pentachlorophenol	16.912	266	1761	0.904	ng/ul#	50
15) Phenanthrene	17.300	178	10904	0.400	ng/ul	97
16) Anthracene	17.397	178	9145	0.398	ng/ul	97
19) Fluoranthene	19.321	202	12935	0.401	ng/ul	100
20) Pyrene	19.679	202	14030	0.424	ng/ul	99
21) Benzo(a)anthracene	21.444	228	9313	0.367	ng/ul	98
22) Chrysene	21.497	228	11787	0.425	ng/ul	99
24) Benzo(b)fluoranthene	23.101	252	10479	0.472	ng/ul	99
25) Benzo(k)fluoranthene	23.148	252	12506	0.508	ng/ul	99
26) Benzo(a)pyrene	23.718	252	9718	0.466	ng/ul	93
27) Indeno(1,2,3-cd)pyrene	26.286	276	10691	0.422	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.326	278	8022	0.404	ng/ul#	89
29) Benzo(g,h,i)perylene	27.017	276	10315	0.449	ng/ul	98

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

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