

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN122221\
 Data File : BN018151.D
 Acq On : 22 Dec 2021 15:23
 Operator : CG/JU
 Sample : PB141448BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS448

Quant Time: Dec 22 23:51:29 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN122121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 22 14:08:04 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	3276	0.400	ng/ul	0.00
4) Naphthalene-d8	10.394	136	10433	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.254	164	6572	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.995	188	14653	0.400	ng/ul	0.00
17) Chrysene-d12	21.191	240	14874	0.400	ng/ul	0.00
23) Perylene-d12	23.421	264	14338	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.076	96	2571	0.759	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.994	152	5312	0.370	ng/ul	0.00
18) Fluoranthene-d10	19.032	212	15588	0.386	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.109	88	7245	1.769	ng/ul#	89
5) Naphthalene	10.443	128	10904	0.368	ng/ul	100
7) 2-Methylnaphthalene	12.071	142	6672	0.350	ng/ul	100
8) 1-Methylnaphthalene	12.286	142	6910	0.349	ng/ul	99
10) Acenaphthylene	13.972	152	9218	0.355	ng/ul	99
11) Acenaphthene	14.314	153	8091	0.353	ng/ul	99
12) Fluorene	15.309	166	9248	0.349	ng/ul	97
14) Pentachlorophenol	16.662	266	1582	0.575	ng/ul	98
15) Phenanthrene	17.038	178	16919	0.354	ng/ul	99
16) Anthracene	17.130	178	13217	0.341	ng/ul	99
19) Fluoranthene	19.060	202	21111	0.370	ng/ul	99
20) Pyrene	19.422	202	22015	0.368	ng/ul	98
21) Benzo(a)anthracene	21.177	228	16536	0.334	ng/ul	99
22) Chrysene	21.229	228	23026	0.380	ng/ul	99
24) Benzo(b)fluoranthene	22.752	252	20266	0.369	ng/ul	99
25) Benzo(k)fluoranthene	22.796	252	26104	0.382	ng/ul	99
26) Benzo(a)pyrene	23.325	252	18751	0.361	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.681	276	24020	0.351	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.701	278	18062	0.348	ng/ul	99
29) Benzo(g,h,i)perylene	26.368	276	20345	0.360	ng/ul	100

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

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