

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102221\
 Data File : BN017039.D
 Acq On : 21 Oct 2021 10:03
 Operator : CG/JU
 Sample : PB140017BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS017

Quant Time: Oct 21 10:39:16 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN101921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 21 10:28:56 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.548	152	1407	0.400	ng/ul	0.00
4) Naphthalene-d8	10.321	136	5273	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.197	164	3072	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.952	188	7070	0.400	ng/ul	0.00
17) Chrysene-d12	21.161	240	7848	0.400	ng/ul	0.00
23) Perylene-d12	23.370	264	9011	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.076	96	938	0.613	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.921	152	2360	0.314	ng/ul	0.00
18) Fluoranthene-d10	18.993	212	6377	0.329	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.109	88	2535	1.378	ng/ul#	45
5) Naphthalene	10.370	128	5124	0.321	ng/ul	98
7) 2-Methylnaphthalene	11.998	142	3233	0.325	ng/ul	100
8) 1-Methylnaphthalene	12.218	142	3307	0.327	ng/ul	100
10) Acenaphthylene	13.915	152	4346	0.350	ng/ul	99
11) Acenaphthene	14.262	153	3724	0.349	ng/ul	98
12) Fluorene	15.257	166	4109	0.333	ng/ul	99
14) Pentachlorophenol	16.610	266	718	0.422	ng/ul	98
15) Phenanthrene	16.994	178	7649	0.341	ng/ul	99
16) Anthracene	17.087	178	6019	0.318	ng/ul	98
19) Fluoranthene	19.021	202	9214	0.343	ng/ul	99
20) Pyrene	19.384	202	9557	0.346	ng/ul	98
21) Benzo(a)anthracene	21.146	228	8205	0.321	ng/ul	100
22) Chrysene	21.199	228	11027	0.379	ng/ul	100
24) Benzo(b)fluoranthene	22.710	252	11834	0.349	ng/ul	99
25) Benzo(k)fluoranthene	22.753	252	13861	0.371	ng/ul	99
26) Benzo(a)pyrene	23.277	252	10237	0.338	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.598	276	16434	0.382	ng/ul	99
28) Dibenzo(a,h)anthracene	25.622	278	12787	0.374	ng/ul	98
29) Benzo(g,h,i)perylene	26.279	276	14025	0.376	ng/ul	99

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

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