

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090221\
 Data File : BN016168.D
 Acq On : 03 Sep 2021 10:26
 Operator : CG/JU
 Sample : PB138650BS
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS650

Quant Time: Sep 03 11:44:18 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN090121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 02 13:32:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	2732	0.400	ng/ul	0.00
4) Naphthalene-d8	10.666	136	11645	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.502	164	6768	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.251	188	13552	0.400	ng/ul	0.00
17) Chrysene-d12	21.441	240	13807	0.400	ng/ul	0.00
23) Perylene-d12	23.835	264	13863	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	2508	0.789	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.256	152	7851	0.479	ng/ul	0.00
18) Fluoranthene-d10	19.281	212	18301	0.494	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.344	88	6383	1.837	ng/ul#	86
5) Naphthalene	10.716	128	12683	0.409	ng/ul	100
7) 2-Methylnaphthalene	12.327	142	8876	0.439	ng/ul	99
8) 1-Methylnaphthalene	12.547	142	8339	0.413	ng/ul	99
10) Acenaphthylene	14.220	152	12585	0.492	ng/ul	99
11) Acenaphthene	14.563	153	8987	0.420	ng/ul	99
12) Fluorene	15.553	166	10393	0.424	ng/ul	99
14) Pentachlorophenol	16.905	266	2497	1.137	ng/ul	99
15) Phenanthrene	17.294	178	16668	0.427	ng/ul	99
16) Anthracene	17.382	178	15993	0.474	ng/ul	98
19) Fluoranthene	19.309	202	20018	0.435	ng/ul	99
20) Pyrene	19.672	202	20676	0.444	ng/ul	96
21) Benzo(a)anthracene	21.424	228	21343	0.505	ng/ul	99
22) Chrysene	21.476	228	21483	0.450	ng/ul	99
24) Benzo(b)fluoranthene	23.104	252	21998	0.422	ng/ul	98
25) Benzo(k)fluoranthene	23.151	252	22434	0.428	ng/ul	99
26) Benzo(a)pyrene	23.727	252	19826	0.448	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.313	276	25416	0.445	ng/ul#	93
28) Dibenzo(a,h)anthracene	26.329	278	21157	0.449	ng/ul	100
29) Benzo(g,h,i)perylene	27.074	276	21447	0.432	ng/ul	99

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

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