

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070823\
 Data File : BN026391.D
 Acq On : 08 Jul 2023 17:58
 Operator : MA/JU
 Sample : PB153848BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS848

Quant Time: Jul 08 18:44:20 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 08 09:20:18 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.930	152	5584	0.400	ng/ul	0.00
4) Naphthalene-d8	10.736	136	19289	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.574	164	9292	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.325	188	17618	0.400	ng/ul	0.00
17) Chrysene-d12	21.519	240	11330	0.400	ng/ul	0.00
23) Perylene-d12	23.978	264	11265	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.306	96	3675	0.584	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.331	152	8615	0.290	ng/ul	0.00
18) Fluoranthene-d10	19.353	212	13372	0.310	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.339	88	9423	1.411	ng/ul#	88
5) Naphthalene	10.785	128	17450	0.316	ng/ul	99
7) 2-Methylnaphthalene	12.402	142	10262	0.292	ng/ul	100
8) 1-Methylnaphthalene	12.617	142	11077	0.304	ng/ul	100
10) Acenaphthylene	14.292	152	16119	0.370	ng/ul	99
11) Acenaphthene	14.635	153	11166	0.314	ng/ul	98
12) Fluorene	15.625	166	13124	0.322	ng/ul	99
14) Pentachlorophenol	16.988	266	2425	0.384	ng/ul	99
15) Phenanthrene	17.368	178	17531	0.301	ng/ul	99
16) Anthracene	17.465	178	14947	0.289	ng/ul	98
19) Fluoranthene	19.386	202	18686	0.309	ng/ul	99
20) Pyrene	19.748	202	20082	0.319	ng/ul	97
21) Benzo(a)anthracene	21.505	228	15205	0.325	ng/ul	99
22) Chrysene	21.554	228	15951	0.328	ng/ul	100
24) Benzo(b)fluoranthene	23.229	252	15951	0.332	ng/ul	93
25) Benzo(k)fluoranthene	23.273	252	16963	0.351	ng/ul	95
26) Benzo(a)pyrene	23.875	252	14203	0.340	ng/ul#	90
27) Indeno(1,2,3-cd)pyrene	26.553	276	17449	0.383	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.563	278	13450	0.390	ng/ul	99
29) Benzo(g,h,i)perylene	27.348	276	15528	0.383	ng/ul	97

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

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