

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071723\
 Data File : BN026541.D
 Acq On : 17 Jul 2023 17:46
 Operator : MA/JU
 Sample : PB153870BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS870

Quant Time: Jul 18 03:05:30 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN071223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 17 17:14:21 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.972	152	1501	0.400	ng/ul	0.02
4) Naphthalene-d8	10.763	136	5092	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.584	164	3388	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.334	188	6860	0.400	ng/ul	0.00
17) Chrysene-d12	21.519	240	5961	0.400	ng/ul	# 0.00
23) Perylene-d12	23.939	264	6013	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.297	96	1535	0.776	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.364	152	2193	0.317	ng/ul	0.02
18) Fluoranthene-d10	19.358	212	6880	0.333	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.331	88	3877	2.073	ng/ul	90
5) Naphthalene	10.813	128	4213	0.301	ng/ul#	93
7) 2-Methylnaphthalene	12.430	142	2426	0.306	ng/ul	99
8) 1-Methylnaphthalene	12.639	142	2835	0.334	ng/ul	100
10) Acenaphthylene	14.311	152	4959	0.299	ng/ul	98
11) Acenaphthene	14.648	153	3813	0.302	ng/ul	97
12) Fluorene	15.643	166	4194	0.316	ng/ul	98
14) Pentachlorophenol	17.000	266	1097	0.634	ng/ul	99
15) Phenanthrene	17.376	178	6490	0.297	ng/ul	99
16) Anthracene	17.473	178	5748	0.303	ng/ul	99
19) Fluoranthene	19.390	202	8525	0.305	ng/ul#	95
20) Pyrene	19.753	202	9206	0.303	ng/ul#	95
21) Benzo(a)anthracene	21.505	228	6735	0.307	ng/ul	99
22) Chrysene	21.560	228	7924	0.318	ng/ul	99
24) Benzo(b)fluoranthene	23.200	252	7284	0.311	ng/ul	95
25) Benzo(k)fluoranthene	23.247	252	7822	0.319	ng/ul#	96
26) Benzo(a)pyrene	23.837	252	6679	0.320	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	26.456	276	9465	0.346	ng/ul#	78
28) Dibenzo(a,h)anthracene	26.469	278	7498	0.357	ng/ul#	90
29) Benzo(g,h,i)perylene	27.234	276	8237	0.341	ng/ul#	89

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

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