

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092823\
 Data File : BN027981.D
 Acq On : 30 Sep 2023 03:55
 Operator : MA/JU
 Sample : PB155863BS
 Misc :
 ALS Vial : 69 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS863

Quant Time: Sep 30 04:24:16 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN091523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 27 03:52:19 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.958	152	7207	0.400	ng/ul	0.00
4) Naphthalene-d8	10.773	136	26778	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.601	164	15602	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.354	188	31723	0.400	ng/ul	0.00
17) Chrysene-d12	21.539	240	20977	0.400	ng/ul	0.00
23) Perylene-d12	24.015	264	20919	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.330	96	4747	0.523	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.357	152	13983	0.345	ng/ul	0.00
18) Fluoranthene-d10	19.380	212	26104	0.407	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.363	88	13478	1.427	ng/ul#	84
5) Naphthalene	10.828	128	21598	0.282	ng/ul	99
7) 2-Methylnaphthalene	12.429	142	14101	0.290	ng/ul	100
8) 1-Methylnaphthalene	12.649	142	13700	0.273	ng/ul	99
10) Acenaphthylene	14.328	152	21768	0.319	ng/ul	99
11) Acenaphthene	14.661	153	15471	0.263	ng/ul	98
12) Fluorene	15.651	166	17002	0.245	ng/ul	99
14) Pentachlorophenol	16.987	266	3530	0.476	ng/ul	99
15) Phenanthrene	17.396	178	26070	0.262	ng/ul	99
16) Anthracene	17.489	178	24350	0.292	ng/ul	99
19) Fluoranthene	19.413	202	27645	0.326	ng/ul	100
20) Pyrene	19.775	202	27921	0.320	ng/ul	99
21) Benzo(a)anthracene	21.522	228	22926	0.282	ng/ul	99
22) Chrysene	21.574	228	21734	0.259	ng/ul	99
24) Benzo(b)fluoranthene	23.243	252	20605	0.221	ng/ul	92
25) Benzo(k)fluoranthene	23.296	252	21192	0.234	ng/ul	92
26) Benzo(a)pyrene	23.901	252	19583	0.261	ng/ul	93
27) Indeno(1,2,3-cd)pyrene	26.620	276	24098	0.251	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.633	278	19730	0.255	ng/ul	95
29) Benzo(g,h,i)perylene	27.431	276	19791	0.243	ng/ul	96

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

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