

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092823\
 Data File : BN027982.D
 Acq On : 30 Sep 2023 04:32
 Operator : MA/JU
 Sample : PB155866BS
 Misc :
 ALS Vial : 70 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS866

Quant Time: Sep 30 05:45:26 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	7512	0.400	ng/ul	0.00
4) Naphthalene-d8	10.779	136	28321	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.601	164	16414	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.354	188	33671	0.400	ng/ul	0.00
17) Chrysene-d12	21.539	240	23592	0.400	ng/ul	0.00
23) Perylene-d12	24.015	264	23343	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.330	96	4646	0.491	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.357	152	13937	0.325	ng/ul	0.00
18) Fluoranthene-d10	19.380	212	26689	0.370	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.363	88	13297	1.350	ng/ul#	83
5) Naphthalene	10.828	128	21482	0.265	ng/ul	100
7) 2-Methylnaphthalene	12.434	142	14025	0.273	ng/ul	100
8) 1-Methylnaphthalene	12.649	142	13667	0.258	ng/ul	99
10) Acenaphthylene	14.328	152	21569	0.301	ng/ul	99
11) Acenaphthene	14.661	153	15318	0.247	ng/ul	98
12) Fluorene	15.651	166	16928	0.232	ng/ul	99
14) Pentachlorophenol	16.991	266	3518	0.446	ng/ul	99
15) Phenanthrene	17.396	178	26184	0.247	ng/ul	99
16) Anthracene	17.489	178	24367	0.275	ng/ul	99
19) Fluoranthene	19.408	202	28488	0.299	ng/ul	97
20) Pyrene	19.775	202	28835	0.293	ng/ul	99
21) Benzo(a)anthracene	21.522	228	24351	0.267	ng/ul	99
22) Chrysene	21.574	228	22900	0.243	ng/ul	99
24) Benzo(b)fluoranthene	23.246	252	21963	0.211	ng/ul	93
25) Benzo(k)fluoranthene	23.296	252	21782	0.216	ng/ul	94
26) Benzo(a)pyrene	23.901	252	20748	0.248	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.619	276	25162	0.235	ng/ul#	92
28) Dibenzo(a,h)anthracene	26.636	278	20668	0.239	ng/ul	95
29) Benzo(g,h,i)perylene	27.434	276	20595	0.226	ng/ul	96

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

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