

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092823\
 Data File : BN027983.D
 Acq On : 30 Sep 2023 05:09
 Operator : MA/JU
 Sample : PB155872BS
 Misc :
 ALS Vial : 71 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS872

Quant Time: Sep 30 05:49:10 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	8187	0.400	ng/ul	0.00
4) Naphthalene-d8	10.773	136	30710	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.597	164	17595	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.354	188	35860	0.400	ng/ul	0.00
17) Chrysene-d12	21.539	240	24175	0.400	ng/ul	0.00
23) Perylene-d12	24.012	264	24178	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.330	96	4482	0.435	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.357	152	13370	0.287	ng/ul	0.00
18) Fluoranthene-d10	19.380	212	25017	0.339	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.364	88	12882	1.200	ng/ul#	83
5) Naphthalene	10.828	128	20598	0.235	ng/ul	100
7) 2-Methylnaphthalene	12.429	142	13477	0.242	ng/ul	100
8) 1-Methylnaphthalene	12.649	142	13095	0.228	ng/ul	98
10) Acenaphthylene	14.328	152	20639	0.269	ng/ul	99
11) Acenaphthene	14.661	153	14759	0.222	ng/ul	97
12) Fluorene	15.647	166	16275	0.208	ng/ul	99
14) Pentachlorophenol	16.991	266	3262	0.389	ng/ul	99
15) Phenanthrene	17.396	178	24996	0.222	ng/ul	99
16) Anthracene	17.489	178	23181	0.246	ng/ul	99
19) Fluoranthene	19.408	202	26589	0.272	ng/ul	97
20) Pyrene	19.775	202	26704	0.265	ng/ul	99
21) Benzo(a)anthracene	21.521	228	22302	0.238	ng/ul	99
22) Chrysene	21.577	228	21078	0.218	ng/ul	99
24) Benzo(b)fluoranthene	23.243	252	20050	0.186	ng/ul	93
25) Benzo(k)fluoranthene	23.293	252	20437	0.196	ng/ul	93
26) Benzo(a)pyrene	23.898	252	18878	0.218	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	26.609	276	23536	0.212	ng/ul#	92
28) Dibenzo(a,h)anthracene	26.632	278	19336	0.216	ng/ul	95
29) Benzo(g,h,i)perylene	27.434	276	19534	0.207	ng/ul	97

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

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