

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100323\
 Data File : BN028014.D
 Acq On : 03 Oct 2023 17:14
 Operator : MA/JU
 Sample : PB155818BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS818

Quant Time: Oct 03 18:05:03 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN100323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 03 15:50:04 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.950	152	4913	0.400	ng/ul	0.00
4) Naphthalene-d8	10.768	136	16398	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.592	164	9567	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.346	188	20411	0.400	ng/ul	0.00
17) Chrysene-d12	21.533	240	19212	0.400	ng/ul	0.00
23) Perylene-d12	24.003	264	19904	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.326	96	3768	0.611	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.352	152	9031	0.342	ng/ul	0.00
18) Fluoranthene-d10	19.371	212	21412	0.351	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.359	88	11370	1.740	ng/ul#	85
5) Naphthalene	10.817	128	14635	0.314	ng/ul	100
7) 2-Methylnaphthalene	12.423	142	9447	0.321	ng/ul	99
8) 1-Methylnaphthalene	12.643	142	9336	0.307	ng/ul	99
10) Acenaphthylene	14.323	152	14313	0.330	ng/ul	100
11) Acenaphthene	14.657	153	10999	0.318	ng/ul	99
12) Fluorene	15.642	166	12262	0.319	ng/ul	99
14) Pentachlorophenol	16.983	266	3102	0.709	ng/ul	100
15) Phenanthrene	17.388	178	20016	0.329	ng/ul	100
16) Anthracene	17.481	178	17938	0.330	ng/ul	100
19) Fluoranthene	19.404	202	23974	0.327	ng/ul	99
20) Pyrene	19.771	202	24773	0.326	ng/ul	98
21) Benzo(a)anthracene	21.516	228	24257	0.346	ng/ul	100
22) Chrysene	21.568	228	24146	0.341	ng/ul	99
24) Benzo(b)fluoranthene	23.237	252	24865	0.337	ng/ul	99
25) Benzo(k)fluoranthene	23.287	252	24924	0.339	ng/ul	99
26) Benzo(a)pyrene	23.892	252	22440	0.357	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.603	276	24666	0.324	ng/ul	98
28) Dibenzo(a,h)anthracene	26.623	278	20101	0.325	ng/ul	99
29) Benzo(g,h,i)perylene	27.414	276	19702	0.316	ng/ul	99

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

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