

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041123\
 Data File : BN024837.D
 Acq On : 11 Apr 2023 15:57
 Operator : CG/JU
 Sample : PB152001BS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS001

Quant Time: Apr 12 01:10:46 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN040723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Apr 08 00:52:14 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.950	152	7372	0.400	ng/u1	0.00
4) Naphthalene-d8	10.759	136	21786	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.593	164	12224	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.338	188	26091	0.400	ng/u1	0.00
17) Chrysene-d12	21.538	240	22220	0.400	ng/u1	0.01
23) Perylene-d12	23.981	264	19388	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.318	96	5726	0.711	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.348	152	11127	0.398	ng/u1	0.00
18) Fluoranthene-d10	19.368	212	23945	0.434	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.352	88	15225	1.717	ng/u1	95
5) Naphthalene	10.814	128	21846	0.393	ng/u1	100
7) 2-Methylnaphthalene	12.425	142	13722	0.397	ng/u1	99
8) 1-Methylnaphthalene	12.640	142	14643	0.400	ng/u1	99
10) Acenaphthylene	14.316	152	18094	0.395	ng/u1	99
11) Acenaphthene	14.658	153	15514	0.395	ng/u1	99
12) Fluorene	15.643	166	17365	0.395	ng/u1	98
14) Pentachlorophenol	16.992	266	2445	0.439	ng/u1	97
15) Phenanthrene	17.380	178	28396	0.386	ng/u1	100
16) Anthracene	17.473	178	23601	0.386	ng/u1	99
19) Fluoranthene	19.400	202	32553	0.425	ng/u1	99
20) Pyrene	19.763	202	33311	0.425	ng/u1	99
21) Benzo(a)anthracene	21.520	228	29453	0.420	ng/u1	100
22) Chrysene	21.575	228	32105	0.423	ng/u1	100
24) Benzo(b)fluoranthene	23.233	252	30948	0.398	ng/u1	98
25) Benzo(k)fluoranthene	23.282	252	30402	0.401	ng/u1	98
26) Benzo(a)pyrene	23.876	252	25161	0.390	ng/u1	96
27) Indeno(1,2,3-cd)pyrene	26.520	276	28161	0.370	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.544	278	21849	0.366	ng/u1	98
29) Benzo(g,h,i)perylene	27.302	276	24738	0.384	ng/u1	99

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

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