

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041323\
 Data File : BN024870.D
 Acq On : 13 Apr 2023 20:57
 Operator : CG/JU
 Sample : PB152109BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS109

Quant Time: Apr 14 02:21:35 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 : Thu Apr 13 02:44:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.951	152	7861	0.400	ng/u1	0.00
4) Naphthalene-d8	10.759	136	23262	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.593	164	12686	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.339	188	27800	0.400	ng/u1	0.00
17) Chrysene-d12	21.529	240	23142	0.400	ng/u1	0.00
23) Perylene-d12	23.970	264	20454	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.318	96	6201	0.722	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.349	152	11033	0.369	ng/u1	0.00
18) Fluoranthene-d10	19.368	212	23909	0.416	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.352	88	16019	1.694	ng/u1	95
5) Naphthalene	10.809	128	21667	0.365	ng/u1	100
7) 2-Methylnaphthalene	12.426	142	13267	0.360	ng/u1	100
8) 1-Methylnaphthalene	12.640	142	14048	0.359	ng/u1	99
10) Acenaphthylene	14.311	152	17175	0.361	ng/u1	99
11) Acenaphthene	14.654	153	14901	0.365	ng/u1	99
12) Fluorene	15.639	166	15969	0.350	ng/u1	100
14) Pentachlorophenol	16.988	266	4320	0.729	ng/u1	99
15) Phenanthrene	17.381	178	28261	0.361	ng/u1	99
16) Anthracene	17.474	178	23314	0.358	ng/u1	99
19) Fluoranthene	19.396	202	31948	0.401	ng/u1	100
20) Pyrene	19.759	202	33256	0.407	ng/u1	100
21) Benzo(a)anthracene	21.511	228	28181	0.386	ng/u1	100
22) Chrysene	21.564	228	31264	0.395	ng/u1	99
24) Benzo(b)fluoranthene	23.221	252	31227	0.381	ng/u1	97
25) Benzo(k)fluoranthene	23.271	252	30581	0.382	ng/u1	98
26) Benzo(a)pyrene	23.862	252	25641	0.377	ng/u1	96
27) Indeno(1,2,3-cd)pyrene	26.507	276	28949	0.361	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.527	278	22510	0.357	ng/u1	98
29) Benzo(g,h,i)perylene	27.285	276	25091	0.370	ng/u1	99

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

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