

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041223\
 Data File : BN024851.D
 Acq On : 12 Apr 2023 17:02
 Operator : CG/JU
 Sample : PB152043BS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS043

Quant Time: Apr 13 02:48:21 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	10035	0.400	ng/u1	0.00
4) Naphthalene-d8	10.765	136	29418	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.593	164	16255	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.338	188	35511	0.400	ng/u1	0.00
17) Chrysene-d12	21.535	240	29311	0.400	ng/u1	0.00
23) Perylene-d12	23.979	264	25947	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.318	96	6305	0.575	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.349	152	11766	0.312	ng/u1	0.00
18) Fluoranthene-d10	19.368	212	24908	0.342	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.352	88	16411	1.359	ng/u1	95
5) Naphthalene	10.814	128	23163	0.309	ng/u1	100
7) 2-Methylnaphthalene	12.426	142	14194	0.304	ng/u1	100
8) 1-Methylnaphthalene	12.640	142	15043	0.304	ng/u1	99
10) Acenaphthylene	14.316	152	18353	0.301	ng/u1	99
11) Acenaphthene	14.653	153	15942	0.305	ng/u1	99
12) Fluorene	15.644	166	17624	0.301	ng/u1	97
14) Pentachlorophenol	16.992	266	4997	0.660	ng/u1	99
15) Phenanthrene	17.381	178	29497	0.295	ng/u1	99
16) Anthracene	17.474	178	24310	0.292	ng/u1	99
19) Fluoranthene	19.396	202	33428	0.331	ng/u1	99
20) Pyrene	19.763	202	34290	0.332	ng/u1	99
21) Benzo(a)anthracene	21.517	228	30566	0.330	ng/u1	100
22) Chrysene	21.573	228	32996	0.329	ng/u1	100
24) Benzo(b)fluoranthene	23.230	252	33537	0.322	ng/u1	98
25) Benzo(k)fluoranthene	23.280	252	32647	0.322	ng/u1	99
26) Benzo(a)pyrene	23.870	252	26987	0.313	ng/u1	96
27) Indeno(1,2,3-cd)pyrene	26.521	276	30001	0.295	ng/u1#	96
28) Dibenzo(a,h)anthracene	26.541	278	23389	0.292	ng/u1	98
29) Benzo(g,h,i)perylene	27.299	276	25856	0.300	ng/u1	99

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

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