

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100323\
 Data File : BM042182.D
 Acq On : 04 Oct 2023 03:30
 Operator : MA/JU
 Sample : PB155904BS
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS904

Quant Time: Oct 04 04:36:49 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM100223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 02 19:28:41 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.933	152	4390	0.400	ng/ul	0.00
4) Naphthalene-d8	10.746	136	10822	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.583	164	4402	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.341	188	8683	0.400	ng/ul	0.00
17) Chrysene-d12	21.512	240	4651	0.400	ng/ul	0.00
23) Perylene-d12	23.915	264	4857	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.360	96	4662	0.797	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.330	152	5879	0.396	ng/ul	0.00
18) Fluoranthene-d10	19.362	212	8699	0.498	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.393	88	14093	2.308	ng/ul#	75
5) Naphthalene	10.796	128	14219	0.394	ng/ul	100
7) 2-Methylnaphthalene	12.407	142	8283	0.383	ng/ul	95
8) 1-Methylnaphthalene	12.621	142	7914	0.363	ng/ul	98
10) Acenaphthylene	14.314	152	12269	0.441	ng/ul	99
11) Acenaphthene	14.647	153	9025	0.429	ng/ul	99
12) Fluorene	15.637	166	10112	0.426	ng/ul	97
14) Pentachlorophenol	16.978	266	3645	0.926	ng/ul	98
15) Phenanthrene	17.383	178	15853	0.443	ng/ul	99
16) Anthracene	17.476	178	13708	0.419	ng/ul	99
19) Fluoranthene	19.394	202	18146	0.514	ng/ul	97
20) Pyrene	19.757	202	18965	0.505	ng/ul#	94
21) Benzo(a)anthracene	21.495	228	13393	0.446	ng/ul	100
22) Chrysene	21.547	228	14193	0.486	ng/ul	99
24) Benzo(b)fluoranthene	23.173	252	13683	0.478	ng/ul	98
25) Benzo(k)fluoranthene	23.222	252	14275	0.506	ng/ul	98
26) Benzo(a)pyrene	23.804	252	12710	0.515	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.408	276	16880	0.482	ng/ul#	88
28) Dibenzo(a,h)anthracene	26.431	278	12513	0.474	ng/ul	96
29) Benzo(g,h,i)perylene	27.176	276	14521	0.497	ng/ul	96

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

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