

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092823\
 Data File : BM042133.D
 Acq On : 29 Sep 2023 02:16
 Operator : MA/JU
 Sample : PB155752BS
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS752

Quant Time: Sep 29 04:42:57 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM092223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 29 04:39:33 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.937	152	2821	0.400	ng/ul	0.00
4) Naphthalene-d8	10.752	136	7189	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.583	164	3647	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.341	188	8174	0.400	ng/ul	0.00
17) Chrysene-d12	21.510	240	5965	0.400	ng/ul	0.00
23) Perylene-d12	23.912	264	6326	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	2722	0.708	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.335	152	3813	0.424	ng/ul	0.00
18) Fluoranthene-d10	19.362	212	8004	0.326	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.398	88	8003	1.979	ng/ul#	87
5) Naphthalene	10.801	128	8688	0.361	ng/ul	99
7) 2-Methylnaphthalene	12.407	142	5397	0.376	ng/ul	99
8) 1-Methylnaphthalene	12.627	142	5297	0.363	ng/ul	100
10) Acenaphthylene	14.315	152	8980	0.367	ng/ul	99
11) Acenaphthene	14.648	153	6533	0.349	ng/ul	99
12) Fluorene	15.638	166	7489	0.355	ng/ul	99
14) Pentachlorophenol	16.970	266	2877	0.665	ng/ul	99
15) Phenanthrene	17.384	178	12967	0.342	ng/ul	99
16) Anthracene	17.476	178	11473	0.350	ng/ul	99
19) Fluoranthene	19.390	202	16297	0.242	ng/ul	100
20) Pyrene	19.757	202	17312	0.261	ng/ul	97
21) Benzo(a)anthracene	21.492	228	14226	0.315	ng/ul	99
22) Chrysene	21.545	228	14537	0.311	ng/ul	99
24) Benzo(b)fluoranthene	23.170	252	14794	0.304	ng/ul	98
25) Benzo(k)fluoranthene	23.219	252	14907	0.315	ng/ul	98
26) Benzo(a)pyrene	23.804	252	13308	0.332	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.401	276	16009	0.291	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.421	278	11885	0.299	ng/ul	97
29) Benzo(g,h,i)perylene	27.172	276	13308	0.288	ng/ul	98

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

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