

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072921\
 Data File : BM031253.D
 Acq On : 29 Jul 2021 17:53
 Operator : CG/JU
 Sample : PB137868BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS868

Quant Time: Jul 29 18:30:23 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 29 13:28:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.611	152	1440	0.400	ng/ul	0.00
4) Naphthalene-d8	10.378	136	5327	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.242	164	3049	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.992	188	5868	0.400	ng/ul	0.00
17) Chrysene-d12	21.180	240	4432	0.400	ng/ul	0.00
23) Perylene-d12	23.350	264	4286	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.203	96	844	0.528	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.970	152	2605	0.290	ng/ul	0.00
18) Fluoranthene-d10	19.023	212	4733	0.320	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.234	88	2303	1.425	ng/ul#	80
5) Naphthalene	10.427	128	4732	0.298	ng/ul	99
7) 2-Methylnaphthalene	12.046	142	3163	0.292	ng/ul	100
8) 1-Methylnaphthalene	12.266	142	2937	0.274	ng/ul	100
10) Acenaphthylene	13.962	152	3887	0.299	ng/ul	100
11) Acenaphthene	14.307	153	3173	0.290	ng/ul	99
12) Fluorene	15.300	166	3549	0.282	ng/ul	99
14) Pentachlorophenol	16.645	266	1094	0.582	ng/ul	100
15) Phenanthrene	17.033	178	5442	0.292	ng/ul	99
16) Anthracene	17.123	178	5080	0.291	ng/ul	99
19) Fluoranthene	19.054	202	5962	0.315	ng/ul	97
20) Pyrene	19.416	202	6038	0.315	ng/ul	98
21) Benzo(a)anthracene	21.166	228	5604	0.313	ng/ul	98
22) Chrysene	21.216	228	5835	0.315	ng/ul	99
24) Benzo(b)fluoranthene	22.701	252	5964	0.309	ng/ul	96
25) Benzo(k)fluoranthene	22.744	252	5845	0.293	ng/ul#	97
26) Benzo(a)pyrene	23.253	252	5175	0.306	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	25.497	276	6884	0.315	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.514	278	5537	0.314	ng/ul	99
29) Benzo(g,h,i)perylene	26.148	276	5872	0.315	ng/ul	97

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

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