

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012022\
 Data File : BM033826.D
 Acq On : 21 Jan 2022 16:31
 Operator : CG/JU
 Sample : N1199-08
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0F14

Quant Time: Jan 22 00:06:17 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM010422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 21 01:34:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.946	152	3064	0.400	ng/ul	0.00
4) Naphthalene-d8	10.761	136	11693	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.583	164	7208	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.316	188	13122	0.400	ng/ul	0.00
17) Chrysene-d12	21.480	240	7243	0.400	ng/ul	# 0.00
23) Perylene-d12	23.878	264	7362	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	14703	4.505	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.348	152	5743	0.361	ng/ul	0.00
18) Fluoranthene-d10	19.338	212	10753	0.479	ng/ul	0.00
Target Compounds						
					Qvalue	
7) 2-Methylnaphthalene	12.420	142	730	0.034	ng/ul	98
8) 1-Methylnaphthalene	12.636	142	2302	0.108	ng/ul	99
11) Acenaphthene	14.644	153	600	0.023	ng/ul	93
12) Fluorene	15.626	166	853	0.027	ng/ul#	97
14) Pentachlorophenol	16.973	266	1007	0.194	ng/ul	96
15) Phenanthrene	17.358	178	2847	0.067	ng/ul	97
19) Fluoranthene	19.369	202	843	0.025	ng/ul#	71
21) Benzo(a)anthracene	21.465	228	612	0.020	ng/ul#	73
24) Benzo(b)fluoranthene	23.144	252	953	0.030	ng/ul#	1
25) Benzo(k)fluoranthene	23.192	252	868	0.027	ng/ul#	1
26) Benzo(a)pyrene	23.771	252	714	0.025	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	26.388	276	2825	0.077	ng/ul#	77
28) Dibenzo(a,h)anthracene	26.407	278	1951	0.067	ng/ul#	59
29) Benzo(g,h,i)perylene	27.156	276	2493	0.079	ng/ul#	79

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

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