

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012022\
 Data File : BM033795.D
 Acq On : 20 Jan 2022 18:59
 Operator : CG/JU
 Sample : N1129-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBM04

Quant Time: Jan 21 01:51:03 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	7.949	152	3650	0.400	ng/ul	0.00
4) Naphthalene-d8	10.760	136	13508	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.587	164	7360	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.319	188	13151	0.400	ng/ul	0.00
17) Chrysene-d12	21.485	240	8394	0.400	ng/ul	# 0.00
23) Perylene-d12	23.885	264	8293	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.303	96	9407	2.419	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.353	152	3175	0.173	ng/ul	0.00
18) Fluoranthene-d10	19.342	212	5678	0.218	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	17.361	178	1870	0.044	ng/ul	97
19) Fluoranthene	19.372	202	2471	0.063	ng/ul#	84
20) Pyrene	19.731	202	2072	0.052	ng/ul#	73
21) Benzo(a)anthracene	21.468	228	1212	0.035	ng/ul#	78
22) Chrysene	21.523	228	1683	0.047	ng/ul#	84
24) Benzo(b)fluoranthene	23.153	252	2722	0.075	ng/ul#	1
25) Benzo(k)fluoranthene	23.202	252	1141	0.032	ng/ul#	1
26) Benzo(a)pyrene	23.778	252	1199	0.038	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	26.390	276	1607	0.039	ng/ul#	15
28) Dibenzo(a,h)anthracene	26.410	278	714	0.022	ng/ul#	1
29) Benzo(g,h,i)perylene	27.168	276	1681	0.048	ng/ul#	48

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

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