

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012022\
 Data File : BM033794.D
 Acq On : 20 Jan 2022 18:24
 Operator : CG/JU
 Sample : N1129-12
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBM05

Quant Time: Jan 21 01:50:02 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	3562	0.400	ng/ul	0.00
4) Naphthalene-d8	10.765	136	13136	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.587	164	7309	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.323	188	13104	0.400	ng/ul	0.00
17) Chrysene-d12	21.485	240	7871	0.400	ng/ul	# 0.00
23) Perylene-d12	23.887	264	7745	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.303	96	11029	2.907	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.353	152	3533	0.198	ng/ul	0.00
18) Fluoranthene-d10	19.342	212	6131	0.251	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	17.361	178	1479	0.035	ng/ul#	92
19) Fluoranthene	19.372	202	1287	0.035	ng/ul#	77
20) Pyrene	19.731	202	1084	0.029	ng/ul#	61
21) Benzo(a)anthracene	21.468	228	810	0.025	ng/ul#	56
22) Chrysene	21.521	228	682	0.020	ng/ul#	58
24) Benzo(b)fluoranthene	23.156	252	1609	0.047	ng/ul#	1
25) Benzo(k)fluoranthene	23.204	252	740	0.022	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	26.393	276	1045	0.027	ng/ul#	56
29) Benzo(g,h,i)perylene	27.168	276	1019	0.031	ng/ul#	22

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

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