

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112221\
 Data File : BM033215.D
 Acq On : 22 Nov 2021 13:45
 Operator : CG/JU
 Sample : M4642-10
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 X3810

Quant Time: Nov 22 14:23:34 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM111921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 19 15:41:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.960	152	1972	0.400	ng/ul	0.00
4) Naphthalene-d8	10.751	136	5531	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.576	164	3164	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.312	188	7047	0.400	ng/ul	0.00
17) Chrysene-d12	21.467	240	7170	0.400	ng/ul	0.00
23) Perylene-d12	23.807	264	7094	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.413	96	783	0.422	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.334	152	2081	0.272	ng/ul	0.00
18) Fluoranthene-d10	19.326	212	6628	0.319	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.805	128	9612	0.561	ng/ul	97
7) 2-Methylnaphthalene	12.411	142	5694	0.513	ng/ul	98
12) Fluorene	15.622	166	8192	0.598	ng/ul	96
14) Pentachlorophenol	16.958	266	2437	1.036	ng/ul	98
15) Phenanthrene	17.354	178	13999	0.548	ng/ul	100
19) Fluoranthene	19.356	202	16910	0.490	ng/ul	98
21) Benzo(a)anthracene	21.453	228	22574	0.882	ng/ul	98
25) Benzo(k)fluoranthene	23.143	252	22372	0.706	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	26.201	276	20586	0.613	ng/ul#	83
29) Benzo(g,h,i)perylene	26.935	276	36906	1.225	ng/ul	93

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

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