

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111921\
 Data File : BM033194.D
 Acq On : 20 Nov 2021 02:04
 Operator : CG/JU
 Sample : M4677-04
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H0AB3

Quant Time: Nov 20 02:53:18 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	3365	0.400	ng/ul	0.00
4) Naphthalene-d8	10.756	136	10497	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.575	164	6890	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.312	188	14630	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.467	240	11672	0.400	ng/ul	# 0.00
23) Perylene-d12	23.807	264	9305	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.413	96	14967	4.730	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.339	152	6044	0.416	ng/ul	0.00
18) Fluoranthene-d10	19.330	212	17979	0.532	ng/ul	0.00
Target Compounds						
					Qvalue	
14) Pentachlorophenol	16.962	266	580	0.119	ng/ul	98
16) Anthracene	17.444	178	1204	0.028	ng/ul#	53
21) Benzo(a)anthracene	21.450	228	1058	0.025	ng/ul#	51
22) Chrysene	21.503	228	1082	0.024	ng/ul#	58
24) Benzo(b)fluoranthene	23.097	252	1202	0.029	ng/ul#	1
25) Benzo(k)fluoranthene	23.143	252	1005	0.024	ng/ul#	1

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

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