

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110921\
 Data File : BM032982.D
 Acq On : 11 Nov 2021 09:31
 Operator : CG/JU
 Sample : M4401-06
 Misc :
 ALS Vial : 78 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG343

Quant Time: Nov 11 10:07:23 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM110921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 09 13:31:56 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.459	152	2914	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.230	136	10833	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.114	164	6619	0.400	ng/ul	-0.01
13) Phenanthrene-d10	16.858	188	13988	0.400	ng/ul	0.00
17) Chrysene-d12	21.047	240	10760	0.400	ng/ul	0.00
23) Perylene-d12	23.154	264	9383	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.857	96	13957	3.751	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	11.831	152	4589	0.298	ng/ul	0.00
18) Fluoranthene-d10	18.889	212	12916	0.396	ng/ul	0.00
Target Compounds						
						Qvalue
14) Pentachlorophenol	16.536	266	74	0.020	ng/ul	88
15) Phenanthrene	16.897	178	1058	0.023	ng/ul#	80

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

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