

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110921\
 Data File : BM032950.D
 Acq On : 10 Nov 2021 13:40
 Operator : CG/JU
 Sample : M4400-01
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BG332

Quant Time: Nov 10 14:37:25 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	7.462	152	1928	0.400	ng/ul	0.00
4) Naphthalene-d8	10.239	136	7112	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.121	164	4348	0.400	ng/ul	# 0.00
13) Phenanthrene-d10	16.862	188	9202	0.400	ng/ul	0.00
17) Chrysene-d12	21.052	240	7653	0.400	ng/ul	0.00
23) Perylene-d12	23.159	264	7435	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.860	96	12098	4.914	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.840	152	3026	0.300	ng/ul	0.00
18) Fluoranthene-d10	18.892	212	9218	0.397	ng/ul	0.00
Target Compounds						
					Qvalue	
14) Pentachlorophenol	16.543	266	77	0.032	ng/ul	93
15) Phenanthrene	16.904	178	1011	0.033	ng/ul#	82
19) Fluoranthene	18.923	202	838	0.023	ng/ul#	67
20) Pyrene	19.285	202	829	0.022	ng/ul#	56

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

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