

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012622\
 Data File : BM033882.D
 Acq On : 26 Jan 2022 22:44
 Operator : CG/JU
 Sample : N1230-18
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0Q18

Quant Time: Jan 27 01:43:35 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM012422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 24 15:20:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.939	152	1832	0.400	ng/ul	0.00
4) Naphthalene-d8	10.751	136	6576	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.576	164	3948	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.309	188	7850	0.400	ng/ul	-0.01
17) Chrysene-d12	21.475	240	6509	0.400	ng/ul #	0.00
23) Perylene-d12	23.868	264	7121	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.296	96	10659	5.566	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.339	152	3443	0.380	ng/ul	0.00
18) Fluoranthene-d10	19.330	212	8315	0.415	ng/ul	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.334	88	258	0.129	ng/ul#	71
12) Fluorene	15.618	166	1292	0.080	ng/ul	94
14) Pentachlorophenol	16.962	266	200783	55.667	ng/ul	97
15) Phenanthrene	17.351	178	946	0.038	ng/ul#	79

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

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