

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012622\
 Data File : BM033873.D
 Acq On : 26 Jan 2022 16:48
 Operator : CG/JU
 Sample : N1230-10
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0Q44

Quant Time: Jan 27 01:42:02 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	1823	0.400	ng/ul	0.00
4) Naphthalene-d8	10.751	136	6765	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.576	164	4569	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.309	188	8176	0.400	ng/ul	-0.01
17) Chrysene-d12	21.477	240	6415	0.400	ng/ul #	0.00
23) Perylene-d12	23.873	264	6773	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.292	96	10473	5.495	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.339	152	4553	0.488	ng/ul	0.00
18) Fluoranthene-d10	19.334	212	10144	0.513	ng/ul	0.00
Target Compounds						Qvalue
5) Naphthalene	10.801	128	13118	0.706	ng/ul#	93
7) 2-Methylnaphthalene	12.411	142	21876	1.811	ng/ul	99
8) 1-Methylnaphthalene	12.631	142	14386	1.214	ng/ul	100
12) Fluorene	15.618	166	1569	0.084	ng/ul	93
14) Pentachlorophenol	16.966	266	177	0.047	ng/ul#	86
15) Phenanthrene	17.351	178	1510	0.058	ng/ul#	86

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

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