

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051322\
 Data File : BM035059.D
 Acq On : 14 May 2022 10:41
 Operator : CG/JU
 Sample : N2603-06
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW487

Quant Time: May 15 01:30:07 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM051222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 12 18:15:39 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.917	152	3913	0.400	ng/ul	0.00
4) Naphthalene-d8	10.715	136	12374	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.552	164	8676	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.292	188	18601	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.473	240	15104	0.400	ng/ul	# 0.00
23) Perylene-d12	23.861	264	13494	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.356	96	36011	7.683	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.305	152	9758	0.494	ng/ul	0.00
18) Fluoranthene-d10	19.322	212	23948	0.462	ng/ul	0.00
Target Compounds						Qvalue
5) Naphthalene	10.765	128	849	0.021	ng/ul#	80
7) 2-Methylnaphthalene	12.382	142	763	0.028	ng/ul	100
8) 1-Methylnaphthalene	12.596	142	839	0.034	ng/ul	93
15) Phenanthrene	17.334	178	2926	0.041	ng/ul#	93
22) Chrysene	21.508	228	2191	0.030	ng/ul#	55
29) Benzo(g,h,i)perylene	27.077	276	1478	0.027	ng/ul#	49

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

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