

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041822\
 Data File : BN019483.D
 Acq On : 18 Apr 2022 17:47
 Operator : CG/JU
 Sample : N2373-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BFFD6

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 04/19/2022
 Supervised By :Yogesh Patel 04/25/2022

Quant Time: Apr 19 03:05:43 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN041822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 18 15:28:08 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.814	152	5243	0.400	ng/ul	0.00
4) Naphthalene-d8	10.613	136	15759	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.457	164	9696	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.197	188	19751	0.400	ng/ul	0.00
17) Chrysene-d12	21.395	240	10788	0.400	ng/ul	0.00
23) Perylene-d12	23.730	264	5693m	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.173	96	28171	4.433	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.213	152	7163	0.339	ng/ul	0.00
18) Fluoranthene-d10	19.231	212	19309	0.439	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.662	128	1090	0.024	ng/ul#	81
15) Phenanthrene	17.239	178	1657	0.025	ng/ul#	83
27) Indeno(1,2,3-cd)pyrene	26.186	276	612	0.026	ng/ul#	84
28) Dibenzo(a,h)anthracene	26.219	278	622	0.039	ng/ul#	9

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

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