

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062422\
 Data File : BN020433.D
 Acq On : 25 Jun 2022 13:49
 Operator : CG/JU
 Sample : N3395-04
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0153

Quant Time: Jun 27 03:42:26 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 24 00:17:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.750	152	3030	0.400	ng/ul	0.00
4) Naphthalene-d8	10.529	136	10846	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.377	164	7045	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.116	188	15791	0.400	ng/ul	0.00
17) Chrysene-d12	21.309	240	13405	0.400	ng/ul	0.00
23) Perylene-d12	23.592	264	12327	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.201	96	17645	5.027	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.124	152	5600	0.315	ng/ul	0.00
18) Fluoranthene-d10	19.151	212	16714	0.378	ng/ul	0.00
Target Compounds						
					Qvalue	
19) Fluoranthene	19.179	202	1977	0.032	ng/ul#	80
21) Benzo(a)anthracene	21.295	228	1894	0.039	ng/ul#	86
22) Chrysene	21.344	228	2009	0.037	ng/ul#	84
24) Benzo(b)fluoranthene	22.902	252	2031	0.036	ng/ul#	29
25) Benzo(k)fluoranthene	22.946	252	2117	0.037	ng/ul#	46
27) Indeno(1,2,3-cd)pyrene	25.933	276	2124	0.041	ng/ul#	93
28) Dibenzo(a,h)anthracene	25.940	278	1672	0.040	ng/ul#	56
29) Benzo(g,h,i)perylene	26.641	276	1712	0.039	ng/ul#	76

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

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