

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091922\
 Data File : BN021806.D
 Acq On : 19 Sep 2022 16:14
 Operator : CG/JU
 Sample : N4595-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CB8G9

Quant Time: Sep 20 00:55:54 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN090822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 20 00:20:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.055	152	5400	0.400	ng/ul	0.00
4) Naphthalene-d8	10.883	136	17578	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.712	164	9917	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.460	188	15380	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.653	240	9161	0.400	ng/ul	# 0.00
23) Perylene-d12	24.164	264	9110	0.400	ng/ul	# 0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.363	96	29266	4.405	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.473	152	10441	0.375	ng/ul	0.00
18) Fluoranthene-d10	19.492	212	13525	0.397	ng/ul	0.00
Target Compounds						
					Qvalue	
10) Acenaphthylene	14.444	152	1616	0.039	ng/ul#	1
15) Phenanthrene	17.502	178	1867	0.038	ng/ul#	1
16) Anthracene	17.595	178	19699	0.489	ng/ul#	85
19) Fluoranthene	19.520	202	2904	0.068	ng/ul#	1
20) Pyrene	19.882	202	3933	0.093	ng/ul#	43
21) Benzo(a)anthracene	21.635	228	5597	0.182	ng/ul#	59
22) Chrysene	21.688	228	1341	0.040	ng/ul#	1
24) Benzo(b)fluoranthene	23.389	252	1633	0.037	ng/ul#	1
25) Benzo(k)fluoranthene	23.439	252	1454	0.034	ng/ul#	1
26) Benzo(a)pyrene	24.050	252	990	0.030	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	26.811	276	1786	0.047	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.831	278	1425	0.045	ng/ul#	1
29) Benzo(g,h,i)perylene	27.626	276	1613	0.046	ng/ul#	18

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

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