

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041522\
 Data File : BN019466.D
 Acq On : 15 Apr 2022 17:14
 Operator : CG/JU
 Sample : N2280-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BFFC3

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 04/18/2022
 Supervised By :Yogesh Patel 04/18/2022

Quant Time: Apr 18 05:03:33 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN041322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 13 14:35:39 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.818	152	4271	0.400	ng/ul	0.00
4) Naphthalene-d8	10.618	136	13123	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.457	164	7917	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.197	188	15987	0.400	ng/ul	0.00
17) Chrysene-d12	21.395	240	10404	0.400	ng/ul	0.00
23) Perylene-d12	23.742	264	5324m	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.173	96	10986	1.999	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.218	152	2571	0.138	ng/ul	0.00
18) Fluoranthene-d10	19.231	212	6752	0.175	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	17.239	178	4894	0.092	ng/ul	96
19) Fluoranthene	19.258	202	6393	0.123	ng/ul	96
20) Pyrene	19.621	202	7995m	0.145	ng/ul	
21) Benzo(a)anthracene	21.383	228	1733m	0.054	ng/ul	
22) Chrysene	21.430	228	3473m	0.083	ng/ul	
24) Benzo(b)fluoranthene	23.037	252	2738m	0.111	ng/ul	
25) Benzo(k)fluoranthene	23.075	252	1274m	0.045	ng/ul	
26) Benzo(a)pyrene	23.637	252	1938	0.092	ng/ul#	64
27) Indeno(1,2,3-cd)pyrene	26.179	276	1154	0.049	ng/ul#	81
29) Benzo(g,h,i)perylene	26.923	276	1262	0.056	ng/ul#	68

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

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