

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071919\
 Data File : BN006923.D
 Acq On : 19 Jul 2019 14:26
 Operator : HP/JU
 Sample : K3764-15
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A52F9

Manual Integrations
 APPROVED

mohammad
 7/22/2019 10:02:47 AM

Quant Time: Jul 19 23:48:06 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN062719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 19 23:29:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	2244	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.36	136	7974	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.24	164	4037	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	8263m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	6498m	0.40	ng/ul	0.02
20) Perylene-d12	23.49	264	5873m	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	1648	0.13	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	4327m	0.18	ng/ul	0.00
Target Compounds						
12) Phenanthrene	17.04	178	2318	0.087	ng/ul	Ovalue 94
15) Fluoranthene	19.09	202	6289m	0.191	ng/ul	
17) Pyrene	19.44	202	5505m	0.185	ng/ul	
18) Benzo(a)anthracene	21.24	228	1869m	0.070	ng/ul	
19) Chrysene	21.30	228	2713	0.091	ng/ul#	93
21) Benzo(b)fluoranthene	22.83	252	3867m	0.161	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	1467m	0.058	ng/ul	
23) Benzo(a)pyrene	23.38	252	2340m	0.099	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.76	276	1904m	0.080	ng/ul	
26) Benzo(g,h,i)perylene	26.43	276	1982m	0.099	ng/ul	

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

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