

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101118\
 Data File : BN003076.D
 Acq On : 10 Oct 2018 22:17
 Operator : MJ/SJ
 Sample : J5157-20
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AE2

Manual Integrations
 APPROVED

Sohil
 10/11/2018 5:45:44 PM

Quant Time: Oct 11 08:05:33 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 11 07:04:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1618	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	6453	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	3408	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	6443m	0.40	ng/ul	-0.01
16) Chrysene-d12	21.23	240	4539m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	4324m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	2876	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	4726m	0.26	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	676	0.041	ng/ul#	79
7) Acenaphthylene	13.98	152	1081	0.069	ng/ul#	86
12) Phenanthrene	17.06	178	848m	0.047	ng/ul	
13) Anthracene	17.15	178	1669m	0.098	ng/ul	
15) Fluoranthene	19.09	202	2979m	0.141	ng/ul	
17) Pyrene	19.45	202	4403m	0.250	ng/ul	
18) Benzo(a)anthracene	21.22	228	2234m	0.153	ng/ul	
19) Chrysene	21.27	228	5036	0.340	ng/ul	98
21) Benzo(b)fluoranthene	22.78	252	6686m	0.429	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	2931m	0.182	ng/ul	
23) Benzo(a)pyrene	23.34	252	4584	0.328	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.67	276	3228	0.201	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.67	278	993	0.077	ng/ul#	60
26) Benzo(g,h,i)perylene	26.35	276	3019	0.223	ng/ul	95

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

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