

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110218\
 Data File : BN003398.D
 Acq On : 02 Nov 2018 09:02
 Operator : MJ/SJ
 Sample : J5701-19
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40K4

Manual Integrations
 APPROVED

Sohil
 11/5/2018 4:09:02 PM

Quant Time: Nov 02 22:21:40 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN110118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 02 23:13:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	9433	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	38616	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	23342	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	55147	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	52901	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	46846	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	7990	0.13	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	22445	0.14	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.27	178	32465	0.180	ng/ul	99
13) Anthracene	17.36	178	6757	0.046	ng/ul	96
15) Fluoranthene	19.29	202	101546	0.500	ng/ul	100
17) Pyrene	19.65	202	86115m	0.448	ng/ul	
18) Benzo(a)anthracene	21.39	228	37229	0.224	ng/ul	98
19) Chrysene	21.45	228	45593	0.242	ng/ul	98
21) Benzo(b)fluoranthene	23.02	252	54461m	0.270	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	21714m	0.111	ng/ul	
23) Benzo(a)pyrene	23.62	252	33486	0.199	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	26.08	276	20808	0.115	ng/ul#	97
25) Dibenzo(a,h)anthracene	26.09	278	5350	0.036	ng/ul#	67
26) Benzo(g,h,i)perylene	26.80	276	16645	0.108	ng/ul	96

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

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