

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101218\
 Data File : BN003113.D
 Acq On : 11 Oct 2018 21:12
 Operator : MJ/SJ
 Sample : J5368-02
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Oct 12 06:51:47 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 : Fri Oct 12 05:29:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1685	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	7290	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	4335	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	9500m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	6810m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	4937m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	3980	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	9750m	0.36	ng/ul	0.00
Target Compounds				Ovalue		
9) Fluorene	15.32	166	354	0.021	ng/ul#	87
12) Phenanthrene	17.06	178	5674	0.214	ng/ul	98
13) Anthracene	17.15	178	1408m	0.056	ng/ul	
15) Fluoranthene	19.09	202	7332m	0.236	ng/ul	
17) Pyrene	19.45	202	5822m	0.220	ng/ul	
18) Benzo(a)anthracene	21.22	228	2641	0.121	ng/ul#	88
19) Chrysene	21.27	228	2384	0.107	ng/ul#	88
21) Benzo(b)fluoranthene	22.78	252	2548m	0.143	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	940m	0.051	ng/ul	
23) Benzo(a)pyrene	23.34	252	1652	0.104	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.68	276	1137	0.062	ng/ul#	49
26) Benzo(g,h,i)perylene	26.35	276	1073	0.069	ng/ul#	50

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

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