

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101318\
 Data File : BN003161.D
 Acq On : 13 Oct 2018 04:18
 Operator : MJ/SJ
 Sample : J5368-06DL 10X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :

Quant Time: Oct 13 06:41:59 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 : Sat Oct 13 05:59:15 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	1236	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.41	136	5358	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.27	164	3253	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	7474	0.40	ng/ul	-0.01
16) Chrysene-d12	21.23	240	6614	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	5837	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.06	152	195	0.02	ng/ul	0.03
14) Fluoranthene-d10	19.06	212	511	0.02	ng/ul	-0.01
Target Compounds				Ovalue		
7) Acenaphthylene	13.99	152	351	0.024	ng/ul#	1
8) Acenaphthene	14.33	153	393	0.035	ng/ul#	90
9) Fluorene	15.33	166	441	0.034	ng/ul#	91
12) Phenanthrene	17.07	178	11968	0.574	ng/ul	100
13) Anthracene	17.17	178	3229	0.164	ng/ul#	91
15) Fluoranthene	19.09	202	25025	1.024	ng/ul	99
17) Pyrene	19.45	202	22972	0.894	ng/ul	97
18) Benzo(a)anthracene	21.21	228	15012	0.705	ng/ul	98
19) Chrysene	21.26	228	14280	0.661	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	23599m	1.123	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	8638m	0.396	ng/ul	
23) Benzo(a)pyrene	23.35	252	15324	0.813	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.66	276	11064	0.511	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.65	278	2740	0.158	ng/ul#	84
26) Benzo(g,h,i)perylene	26.34	276	9130	0.500	ng/ul	99

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

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