

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN110118\
 Data File : BN003381.D
 Acq On : 01 Nov 2018 16:38
 Operator : MJ/SJ
 Sample : J5701-19
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40K4

Quant Time: Nov 01 17:52:22 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 : Thu Nov 01 13:18:49 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	9446	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	38081	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	22511	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	53444	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	52434	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	48328	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	7738	0.13	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	21248	0.14	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	17.27	178	30864	0.176	ng/ul	99
13) Anthracene	17.36	178	6013	0.042	ng/ul	95
15) Fluoranthene	19.29	202	96473	0.490	ng/ul	99
17) Pyrene	19.65	202	87960	0.461	ng/ul	99
18) Benzo(a)anthracene	21.40	228	34641	0.211	ng/ul	99
19) Chrysene	21.45	228	44472	0.238	ng/ul	98
21) Benzo(b)fluoranthene	23.03	252	77900	0.374	ng/ul	94
22) Benzo(k)fluoranthene	23.03	252	77900	0.386	ng/ul#	94
23) Benzo(a)pyrene	23.63	252	34359	0.198	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	26.08	276	20966	0.113	ng/ul#	95
25) Dibenzo(a,h)anthracene	26.09	278	5018	0.033	ng/ul#	69
26) Benzo(g,h,i)perylene	26.81	276	15902	0.100	ng/ul	96

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

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