

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110518\
 Data File : BN003428.D
 Acq On : 05 Nov 2018 14:16
 Operator : MJ/SJ
 Sample : J5704-10
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40K2

Quant Time: Nov 05 15:11:09 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 : Mon Nov 05 13:38:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	10631	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	44507	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	25969	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.33	188	2967382	0.40	ng/ul	0.10
16) Chrysene-d12	21.54	240	437	0.40	ng/ul	0.13
20) Perylene-d12	23.87	264	446	0.40	ng/ul	0.15
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	21794	0.31	ng/ul	0.00
14) Fluoranthene-d10	0.00	212	0	0.00	ng/ul	
Target Compounds				Qvalue		
3) Naphthalene	10.70	128	5755	0.046	ng/ul#	94
5) 2-Methylnaphthalene	12.31	142	3191	0.036	ng/ul	99
7) Acenaphthylene	14.21	152	45335	0.402	ng/ul	99
8) Acenaphthene	14.55	153	3924	0.039	ng/ul	99
9) Fluorene	15.54	166	15629	0.132	ng/ul	99
17) Pyrene	19.65	202	721761	454.053	ng/ul	100
18) Benzo(a)anthracene	21.45	228	379127	276.715	ng/ul	94
19) Chrysene	21.56	228	35768	22.970	ng/ul#	60
21) Benzo(b)fluoranthene	23.20	252	65535	34.101	ng/ul	84
22) Benzo(k)fluoranthene	23.20	252	65535	35.208	ng/ul#	84
23) Benzo(a)pyrene	23.77	252	69701	43.432	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	26.35	276	9210	5.361	ng/ul#	50
25) Dibenzo(a,h)anthracene	26.20	278	18399	13.140	ng/ul#	72
26) Benzo(g,h,i)perylene	26.80	276	152941	103.814	ng/ul	98

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

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