

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110118\
 Data File : BN003378.D
 Acq On : 01 Nov 2018 14:51
 Operator : MJ/SJ
 Sample : J5704-13
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40D8

Quant Time: Nov 01 17:25:42 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	6786	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	27444	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	15980	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	40497	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	39254	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	31223	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	13390	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	45567	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.70	128	20878	0.271	ng/ul	99
5) 2-Methylnaphthalene	12.31	142	14806	0.267	ng/ul	100
7) Acenaphthylene	14.21	152	46111	0.664	ng/ul	100
9) Fluorene	15.54	166	25084	0.343	ng/ul	99
12) Phenanthrene	17.28	178	89523	0.674	ng/ul	99
13) Anthracene	17.36	178	27393	0.253	ng/ul	99
15) Fluoranthene	19.29	202	74689	0.500	ng/ul	87
17) Pyrene	19.65	202	84162	0.589	ng/ul	99
18) Benzo(a)anthracene	21.40	228	102304	0.831	ng/ul	100
19) Chrysene	21.45	228	58211	0.416	ng/ul	99
21) Benzo(b)fluoranthene	23.03	252	44687	0.332	ng/ul	95
22) Benzo(k)fluoranthene	23.07	252	69553	0.534	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.09	276	48287	0.402	ng/ul#	98
25) Dibenzo(a,h)anthracene	26.10	278	44107	0.450	ng/ul	97
26) Benzo(g,h,i)perylene	26.81	276	62858	0.609	ng/ul	99

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

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