

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112018\
 Data File : BN003625.D
 Acq On : 20 Nov 2018 10:30
 Operator : MJ/SJ
 Sample : J6002-03
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X2E78

Quant Time: Nov 20 11:13:17 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 20 03:27:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	5200	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	20448	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	11158	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	26985	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	29231	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	24488	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	11962	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	29593	0.37	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.69	128	15028	0.262	ng/ul	99
5) 2-Methylnaphthalene	12.30	142	11039	0.275	ng/ul	100
7) Acenaphthylene	14.20	152	40737	0.811	ng/ul	100
8) Acenaphthene	14.54	153	14635	0.332	ng/ul	98
9) Fluorene	15.53	166	18410	0.359	ng/ul	100
11) Pentachlorophenol	16.87	266	1165	0.240	ng/ul	97
12) Phenanthrene	17.26	178	64949	0.733	ng/ul	100
13) Anthracene	17.36	178	17969	0.247	ng/ul	100
15) Fluoranthene	19.28	202	46045	0.440	ng/ul	99
17) Pyrene	19.64	202	49178	0.469	ng/ul	99
18) Benzo(a)anthracene	21.38	228	64895	0.703	ng/ul	100
19) Chrysene	21.43	228	40360	0.382	ng/ul	100
21) Benzo(b)fluoranthene	23.01	252	36377	0.359	ng/ul	99
22) Benzo(k)fluoranthene	23.01	252	36348	0.368	ng/ul	99
23) Benzo(a)pyrene	23.61	252	43240	0.493	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.06	276	35198	0.350	ng/ul	99
25) Dibenzo(a,h)anthracene	26.08	278	30264	0.374	ng/ul	99
26) Benzo(g,h,i)perylene	26.79	276	35634	0.408	ng/ul	99

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

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