

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110218\
 Data File : BN003416.D
 Acq On : 02 Nov 2018 22:47
 Operator : MJ/SJ
 Sample : J5701-17
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40J3

Manual Integrations
 APPROVED

Sohil
 11/5/2018 4:09:24 PM

Quant Time: Nov 03 00:19:42 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 : Fri Nov 02 23:49:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.86	152	10853	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	45007	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	25553	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	58062m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	45920m	0.40	ng/ul	0.00
20) Perylene-d12	23.73	264	38105m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	14979	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	36131m	0.22	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.70	128	4363	0.035	ng/ul#	93
5) 2-Methylnaphthalene	12.31	142	2081	0.023	ng/ul	100
7) Acenaphthylene	14.21	152	5746	0.052	ng/ul#	95
9) Fluorene	15.54	166	2382	0.020	ng/ul#	97
12) Phenanthrene	17.28	178	40815m	0.214	ng/ul	
13) Anthracene	17.36	178	8759m	0.056	ng/ul	
15) Fluoranthene	19.29	202	83730m	0.391	ng/ul	
17) Pyrene	19.65	202	84831m	0.508	ng/ul	
18) Benzo(a)anthracene	21.39	228	41249m	0.287	ng/ul	
19) Chrysene	21.45	228	42120m	0.257	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	41845m	0.255	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	14474m	0.091	ng/ul	
23) Benzo(a)pyrene	23.63	252	31695m	0.231	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.08	276	20835m	0.142	ng/ul	
25) Dibenzo(a,h)anthracene	26.08	278	6061m	0.051	ng/ul	
26) Benzo(a,h,i)perylene	26.81	276	20640m	0.164	ng/ul	

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

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