

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110218\
 Data File : BN003413.D
 Acq On : 02 Nov 2018 21:01
 Operator : MJ/SJ
 Sample : J5704-21
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A40H9

Manual Integrations
 APPROVED

Sohil
 11/5/2018 4:09:21 PM

Quant Time: Nov 03 00:04:34 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	9777	0.40	ng/ul	0.00
2) Naphthalene-d8	10.65	136	40996	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	24441	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	56795m	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	51950m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	40208m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.24	152	18094	0.28	ng/ul	0.00
14) Fluoranthene-d10	19.26	212	47760m	0.30	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.70	128	3652	0.032	ng/ul#	92
7) Acenaphthylene	14.21	152	17863	0.168	ng/ul	98
8) Acenaphthene	14.55	153	2582	0.027	ng/ul	98
9) Fluorene	15.54	166	7751	0.069	ng/ul#	96
11) Pentachlorophenol	16.87	266	469m	0.041	ng/ul	
12) Phenanthrene	17.28	178	239873m	1.288	ng/ul	
13) Anthracene	17.36	178	14784m	0.097	ng/ul	
15) Fluoranthene	19.28	202	382186m	1.826	ng/ul	
17) Pyrene	19.65	202	289741	1.533	ng/ul	99
18) Benzo(a)anthracene	21.39	228	93234m	0.572	ng/ul	
19) Chrysene	21.44	228	146328m	0.790	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	180117m	1.040	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	68805m	0.410	ng/ul	
23) Benzo(a)pyrene	23.63	252	94676m	0.654	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.08	276	73909m	0.477	ng/ul	
25) Dibenzo(a,h)anthracene	26.09	278	18159m	0.144	ng/ul	
26) Benzo(a,h,i)perylene	26.81	276	49693m	0.374	ng/ul	

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

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