

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121118\
 Data File : BN003825.D
 Acq On : 11 Dec 2018 21:19
 Operator : JU/SJ
 Sample : J6170-19
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9M14

Manual Integrations
 APPROVED

Sohil
 12/12/2018 5:50:29 PM

Quant Time: Dec 12 03:36:37 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 : Wed Dec 12 02:52:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	4935	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	18933	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	8911	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	16682m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	12464m	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	12430m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	5331	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	8034m	0.16	ng/ul	0.00
Target Compounds				Ovalue		
9) Fluorene	15.52	166	1058	0.026	ng/ul#	96
12) Phenanthrene	17.25	178	24586	0.449	ng/ul	99
13) Anthracene	17.34	178	3423m	0.076	ng/ul	
15) Fluoranthene	19.27	202	87607m	1.355	ng/ul	
17) Pyrene	19.63	202	64274m	1.438	ng/ul	
18) Benzo(a)anthracene	21.38	228	34404	0.874	ng/ul	97
19) Chrysene	21.43	228	41263	0.916	ng/ul	99
21) Benzo(b)fluoranthene	23.01	252	61315m	1.192	ng/ul	
22) Benzo(k)fluoranthene	23.05	252	20602m	0.411	ng/ul	
23) Benzo(a)pyrene	23.61	252	33947	0.762	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.06	276	30865	0.604	ng/ul	95
25) Dibenzo(a,h)anthracene	26.07	278	7653	0.186	ng/ul#	88
26) Benzo(g,h,i)perylene	26.78	276	24541	0.553	ng/ul	96

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

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